Sample ID:	6507-1	Operator:	Jeremy Armstrong
Instrument ID:	Varian GC/MS #1	Last Calibration:	4/17/2012 10:51 AM
Acquisition Date:	4/24/2012 3:25 PM	Data File:	...il\042312\6507-1.sms
Calculation Date:	4/24/2012 3:50 PM	Method:	..._cain_041612_jwa.mth
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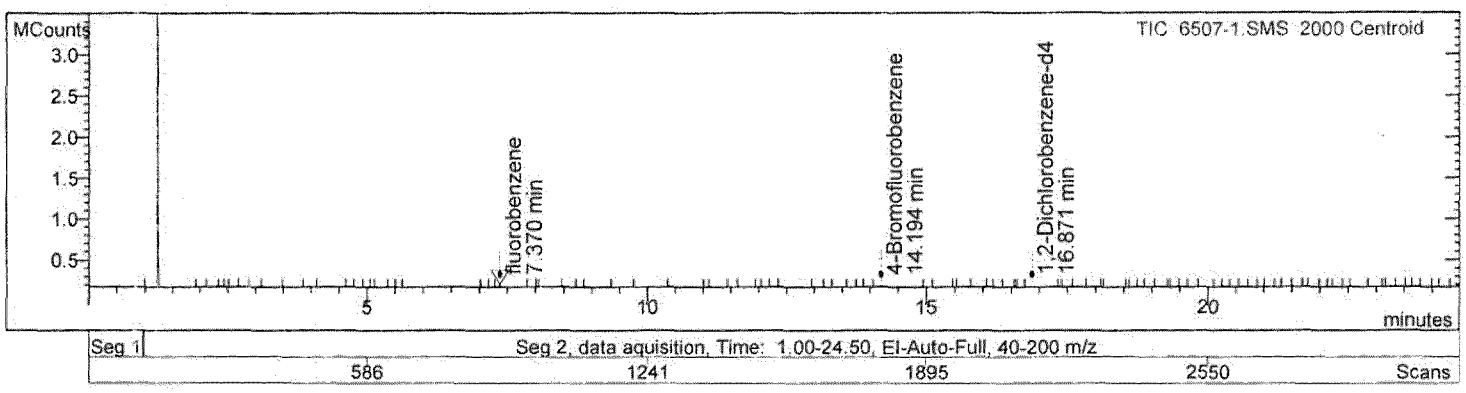
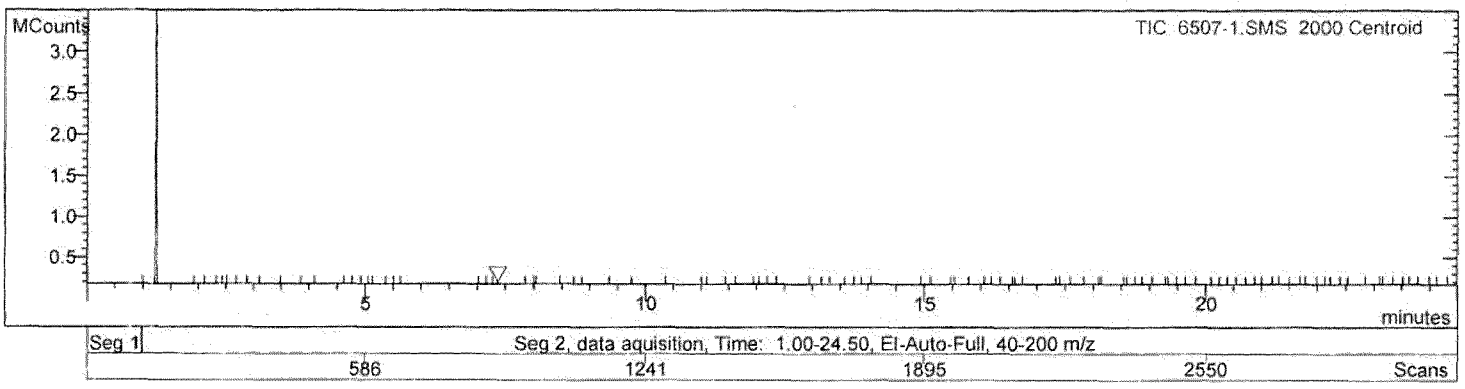
Multiplier:	1.000000		



Target Compounds

#	RT	Peak Name	Res Type	Quan Ions	Area	Amount/RF
1	7.370	fluorobenzene	Id.	96.0	225289	100.000 ug/L
2	3.971	TBA	Miss.	59.0+57.0	0	0.000 ug/L
3	4.166	MTBE	Miss.	73.0+57.0	0	0.000 ug/L
4	4.932	di-isopropyl ether (DIPE)	Miss.	45.0+69.0	0	0.000 ug/L
5	5.470	ethyl tert-butyl ether	Miss.	59.0+87.0	0	0.000 ug/L
6	6.907	benzene	Miss.	78.0+77.0	0	0.000 ug/L
7	7.137	tert-amyl methyl ether (T	Miss.	73.0+87.0	0	0.000 ug/L
8	8.057	2,5-dimethylfuran	Miss.	96.0+95.0	0	0.000 ug/L
9	9.908	toluene	Miss.	91.0+92.0	0	0.000 ug/L
10	11.342	1,2-Dibromoethane	Miss.	108.9	0	0.000 ug/L
11	12.413	ethyl benzene	Miss.	91.0+106.0	0	0.000 ug/L
12	12.622	m+p xylene	Miss.	91.0+106.0	0	0.000 ug/L
13	13.281	o-xylene	Miss.	91.0+106.0	0	0.000 ug/L
14	14.194	4-Bromofluorobenzene	Id.	95.0+176.0	143611	100.633 ug/L
15	14.990	1,3,5-Trimethylbenzene	Miss.	105.0+120.0	0	0.000 ug/L
16	15.634	1,2,4-Trimethylbenzene	Miss.	105.0+120.0	0	0.000 ug/L
17	16.363	1,2,3-Trimethylbenzene	Miss.	105.0+120.0	0	0.000 ug/L
18	16.871	1,2-Dichlorobenzene-d4	Id.	115.0	34059	98.094 ug/L
19	19.901	Naphthalene	Miss.	128.0+127.0	0	0.000 ug/L



Unidentified Peaks

None

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Revision Log

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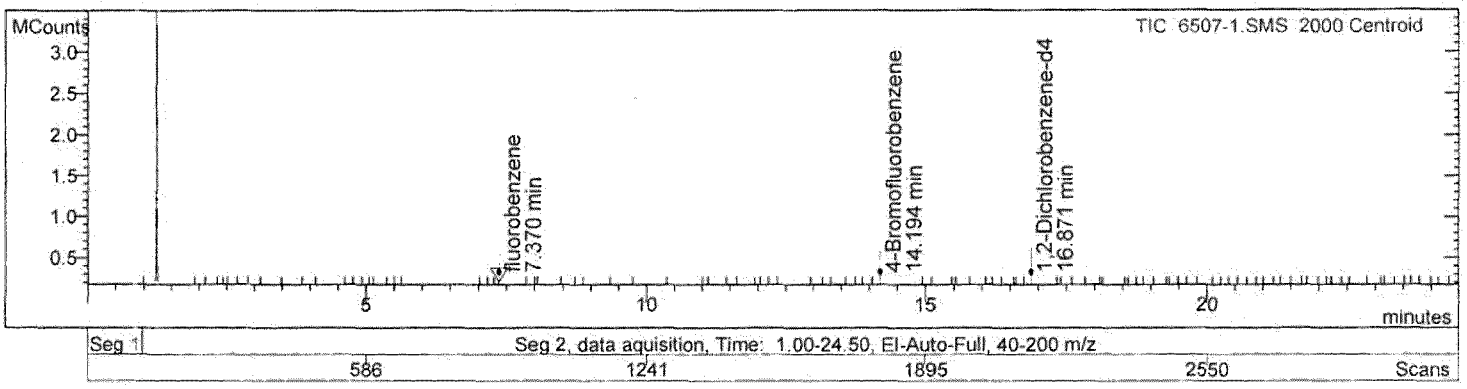
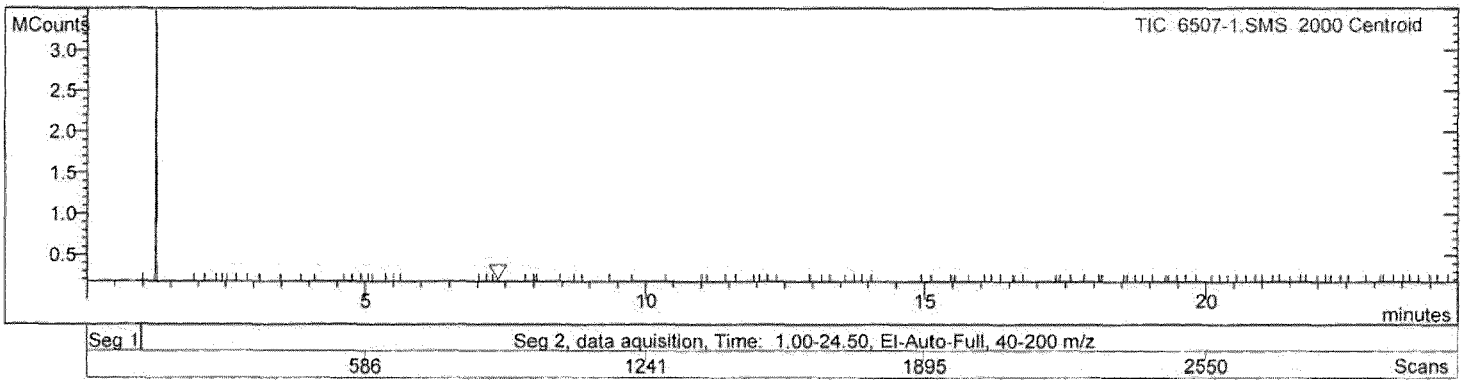
Sample ID:	6507-1	Operator:	Jeremy Armstrong
Instrument ID:	Varian GC/MS #1	Last Calibration:	4/9/2012 8:17 AM
Acquisition Date:	4/24/2012 3:25 PM	Data File:	...il\042312\6507-1.sms
Calculation Date:	4/24/2012 7:18 PM	Method:	..._cain_040612_jwa.mth
Inj. Sample Notes:	FieldBlk 1		
Multiplier:	1.000000		



Target Compounds

#	RT	Peak Name	Res Type	Quan Ions	Area	Amount/RF
1	7.370	fluorobenzene	Id.	96.0	225289	100.000 ug/L
2	2.685	Ethanol	Miss.	45.0+44.0	0	0.000 ug/L
3	3.383	isopropanol	Miss.	45.0+43.0	0	0.000 ug/L
4	5.039	n-propanol	Miss.	59.0+42.0	0	0.000 ug/L
5	6.927	iso-butanol	Miss.	43.1+56.0	0	0.000 ug/L
6	7.939	n-butanol	Miss.	56.0+55.0	0	0.000 ug/L
7	14.194	4-Bromofluorobenzene	Id.	95.0+176.0	143611	101.955 ug/L
8	16.871	1,2-Dichlorobenzene-d4	Id.	115.0	34059	94.983 ug/L



Unidentified Peaks

None

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Revision Log

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Sample ID:	6507-1	Operator:	Jeremy Armstrong
Instrument ID:	Varian GC/MS #1	Last Calibration:	4/16/2012 10:56 AM
Acquisition Date:	4/24/2012 3:25 PM	Data File:	...il\042312\6507-1.sms
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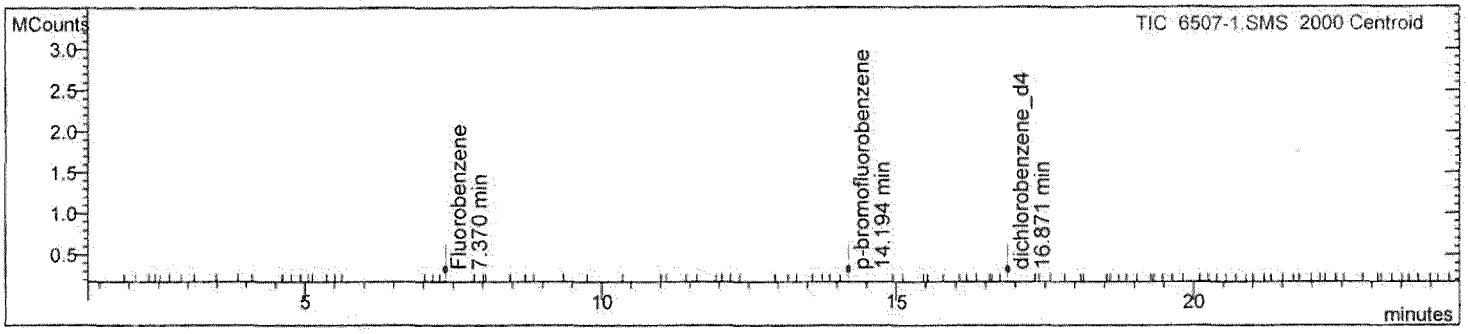
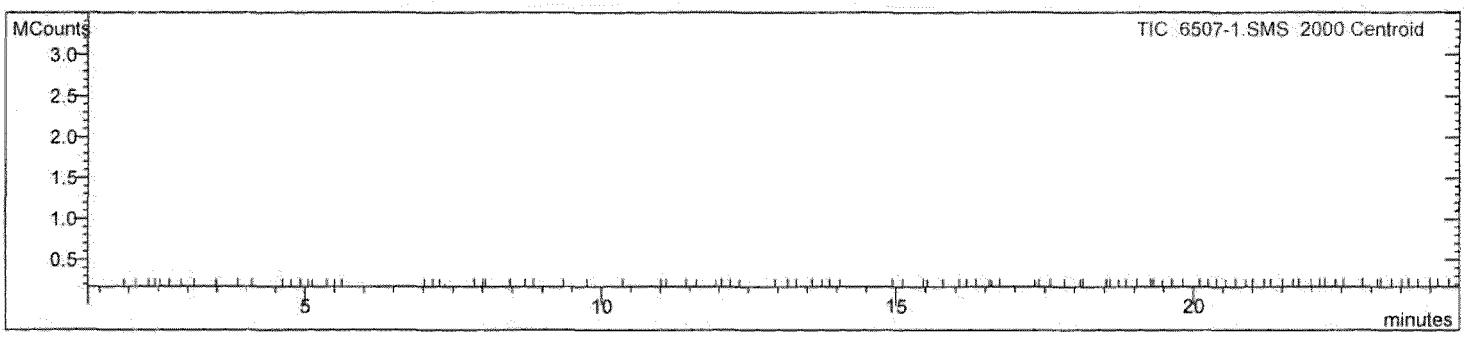
Multiplier:	1.000000		



Target Compounds

#	RT	Peak Name	Res Type	Quan Ions	Area	Amount/RF
1	7.370	Fluorobenzene	Id.	96.0	224812	100.000 ug/L
2	3.160	acetone	Miss.	43.0	0	0.000 ug/L
3	4.000	TBA	Miss.	59.0	0	0.000 ug/L
4	4.194	MTBE	Miss.	73.0	0	0.000 ug/L
5	14.194	p-bromofluorobenzene	Id.	174.0	49546	101.584 ug/L
6	16.871	dichlorobenzene_d4	Id.	115.0	33507	103.685 ug/L



Unidentified Peaks

None

Sample ID:	6507-1	Operator:	Jeremy Armstrong
Instrument ID:	Varian GC/MS #1	Last Calibration:	4/20/2012 12:37 PM
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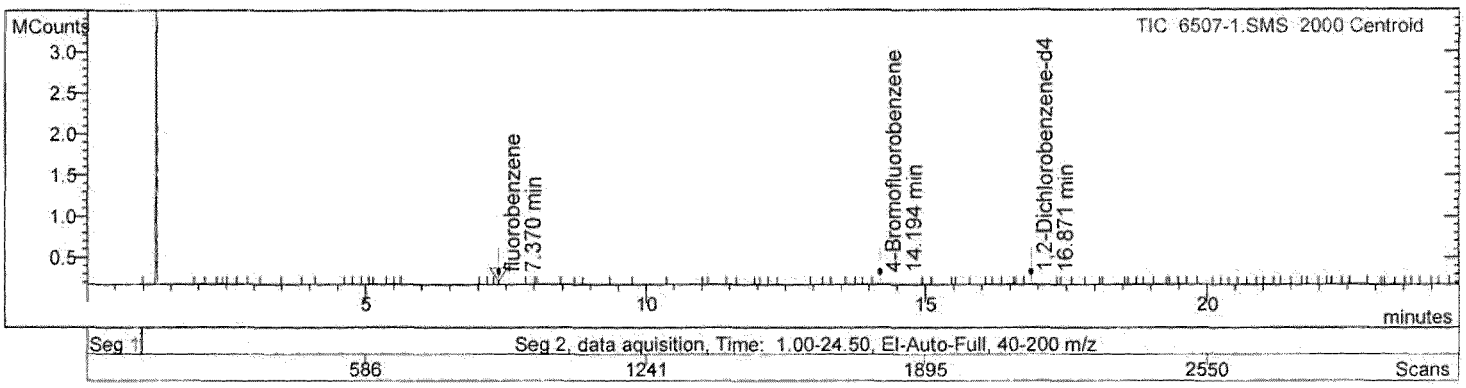
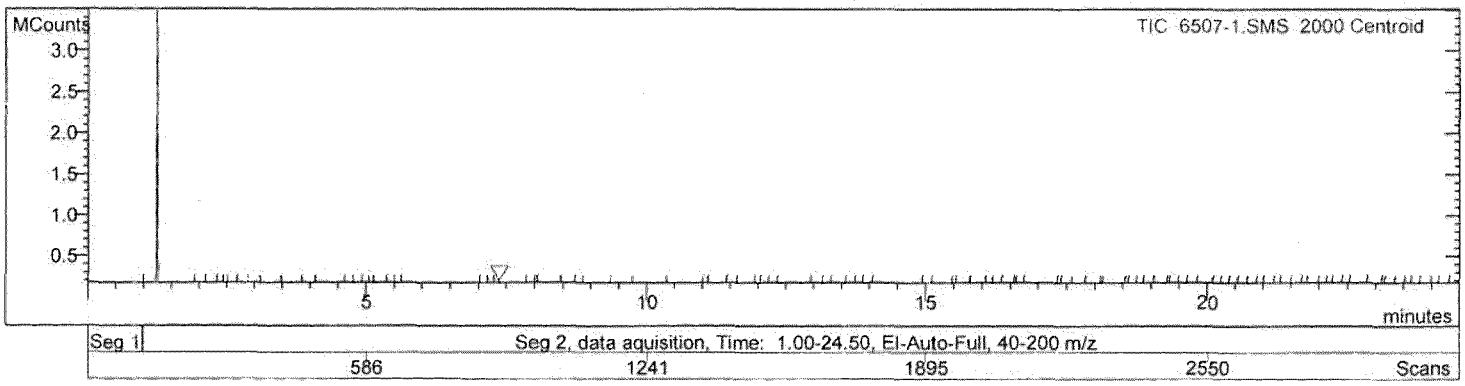


Target Compounds

#	RT	Peak Name	Res Type	Quan Ions	Area	Amount/RF
1	7.370	fluorobenzene	Id.	96.0	225289	100.000 ug/L
2	1.677	Vinyl Chloride	Miss.	62.0+64.0	0	0.000 ug/L
3	3.075	1,1-Dichloroethene	Miss.	61.0+96.0	0	0.000 ug/L
4	3.770	Methylene Chloride	Miss.	49.0+84.0	0	0.000 ug/L
5	4.176	trans-1,2-dichloroethene	Miss.	61.0+96.0	0	0.000 ug/L
6	4.793	1,1-dichloroethane	Miss.	63.0+65.0	0	0.000 ug/L
7	5.661	cis-1,2-dichloroethene	Miss.	61.0+96.0	0	0.000 ug/L
8	6.156	chloroform	Miss.	83.0+85.0	0	0.000 ug/L
9	6.391	1,1,1-trichloroethane	Miss.	97.0+61.0	0	0.000 ug/L
10	6.634	carbon tetrachloride	Miss.	117.0+119.0	0	0.000 ug/L
11	6.980	1,2-dichloroethane	Miss.	62.0+64.0	0	0.000 ug/L
12	7.939	trichloroethene	Miss.	95.0+132.0	0	0.000 ug/L
13	10.613	1,1,2-Trichloroethane	Miss.	96.8	0	0.000 ug/L
14	10.817	tetrachloroethene	Miss.	166.0+129.0	0	0.000 ug/L
15	12.225	chlorobenzene	Miss.	112.0+77.0	0	0.000 ug/L
16	14.194	4-Bromofluorobenzene	Id.	95.0+176.0	143611	102.946 ug/L
17	16.106	1,3-dichlorobenzene	Miss.	146.1+111.0	0	0.000 ug/L
18	16.271	1,4-dichlorobenzene	Miss.	146.1+111.0	0	0.000 ug/L
19	16.871	1,2-Dichlorobenzene-d4	Id.	115.0	34059	106.771 ug/L
20	16.916	1,2-dichlorobenzene	Miss.	146.1	0	0.000 ug/L



Unidentified Peaks

None

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Revision Log

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Print Date: 24 Apr 2012 19:13:19

Chromatogram Plot

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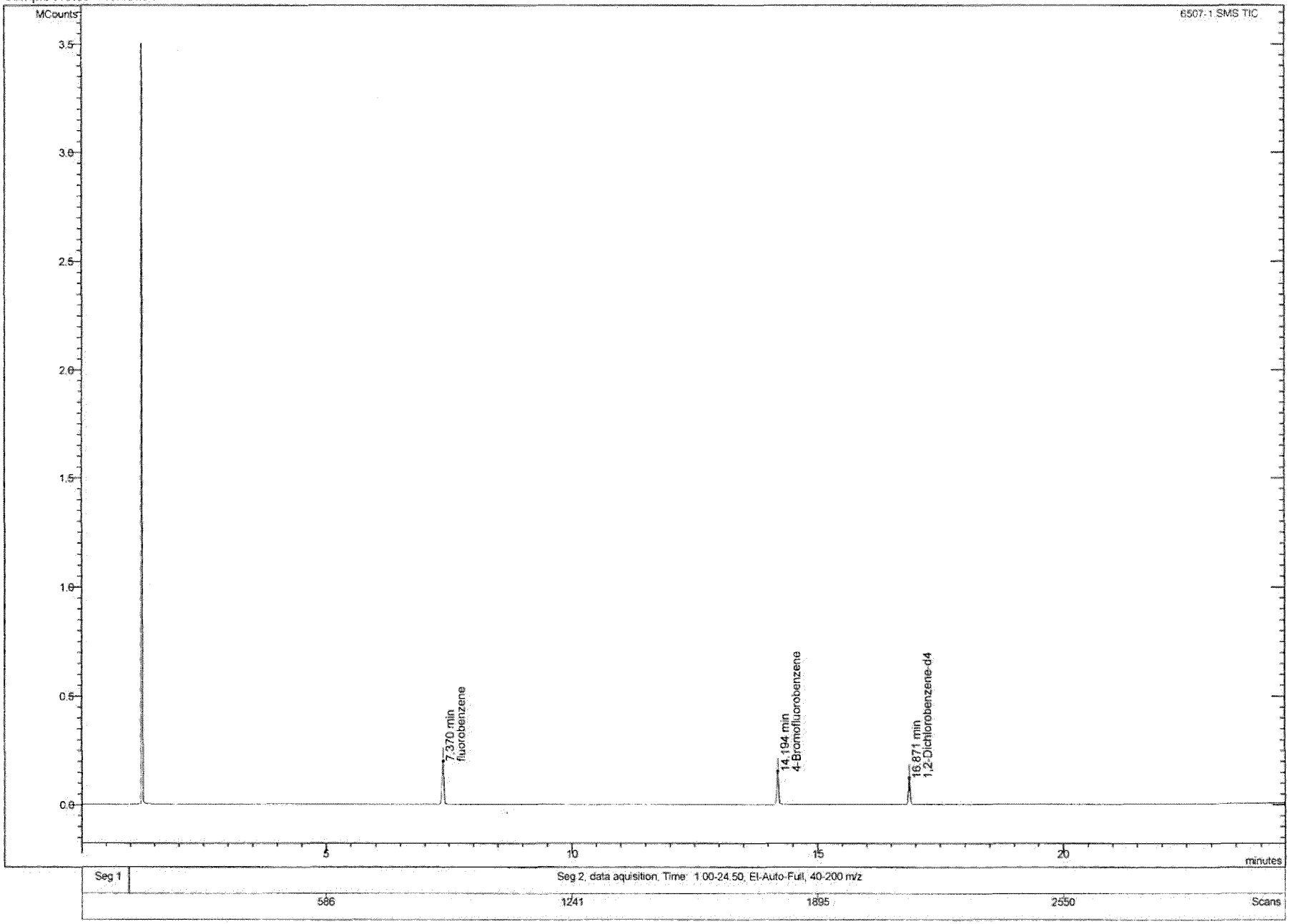
Scan Range: 1 - 3129 Time Range: 0.00 - 24.48 min.
Sample Notes: FieldBlk 1

Operator: Jeremy Armstrong
Date: 4/24/2012 3:25 PM



EPAPAV0043725

Sample ID:	6507-2	Operator:	Jeremy Armstrong
Instrument ID:	Varian GC/MS #1	Last Calibration:	4/17/2012 10:51 AM
Acquisition Date:	4/24/2012 4:16 PM	Data File:	...il\042312\6507-2.sms
Calculation Date:	4/24/2012 4:40 PM	Method:	..._cain_041612_jwa.mth
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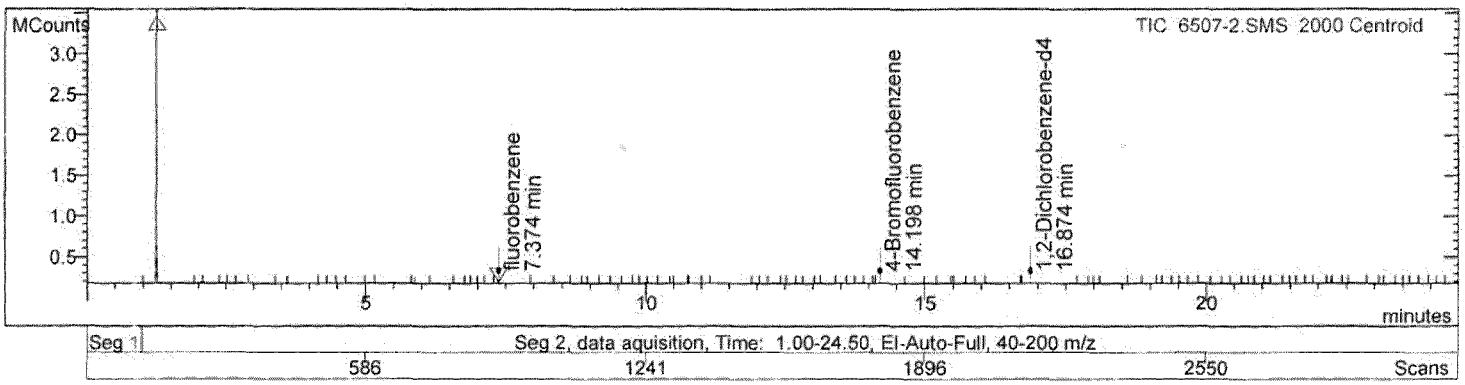
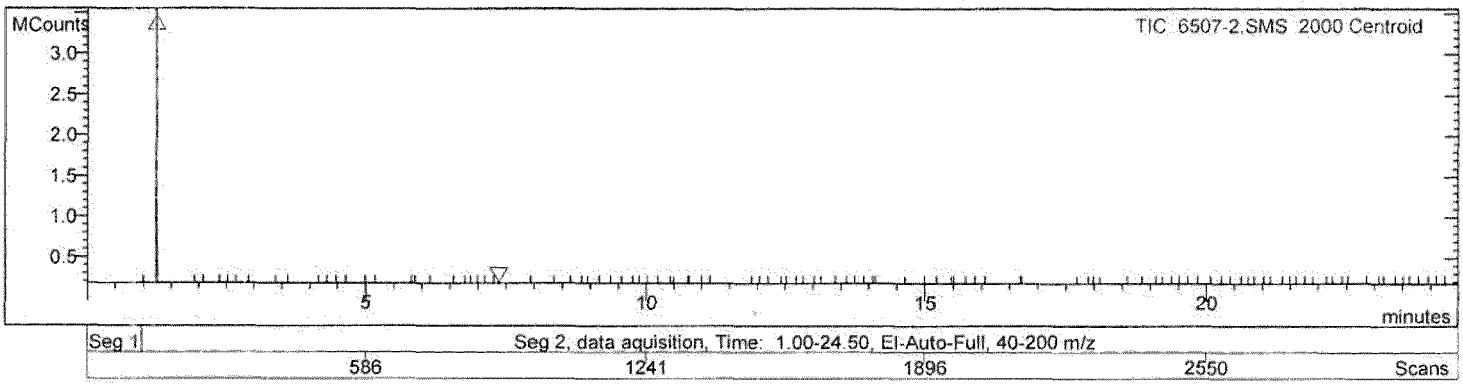


Target Compounds

#	RT	Peak Name	Res Type	Quan Ions	Area	Amount/RF
1	7.374	fluorobenzene	Id.	96.0	212209	100.000 ug/L
2	3.971	TBA	Miss.	59.0+57.0	0	0.000 ug/L
3	4.166	MTBE	Miss.	73.0+57.0	0	0.000 ug/L
4	4.932	di-isopropyl ether (DIPE)	Miss.	45.0+69.0	0	0.000 ug/L
5	5.470	ethyl tert-butyl ether	Miss.	59.0+87.0	0	0.000 ug/L
6	6.907	benzene	Miss.	78.0+77.0	0	0.000 ug/L
7	7.137	tert-amyl methyl ether (T	Miss.	73.0+87.0	0	0.000 ug/L
8	8.057	2,5-dimethylfuran	Miss.	96.0+95.0	0	0.000 ug/L
9	9.908	toluene	Miss.	91.0+92.0	0	0.000 ug/L
10	11.342	1,2-Dibromoethane	Miss.	108.9	0	0.000 ug/L
11	12.413	ethyl benzene	Miss.	91.0+106.0	0	0.000 ug/L
12	12.622	m+p xylene	Miss.	91.0+106.0	0	0.000 ug/L
13	13.281	o-xylene	Miss.	91.0+106.0	0	0.000 ug/L
14	14.198	4-Bromofluorobenzene	Id.	95.0+176.0	138084	102.928 ug/L
15	14.990	1,3,5-Trimethylbenzene	Miss.	105.0+120.0	0	0.000 ug/L
16	15.634	1,2,4-Trimethylbenzene	Miss.	105.0+120.0	0	0.000 ug/L
17	16.363	1,2,3-Trimethylbenzene	Miss.	105.0+120.0	0	0.000 ug/L
18	16.874	1,2-Dichlorobenzene-d4	Id.	115.0	32341	98.889 ug/L
19	19.901	Naphthalene	Miss.	128.0+127.0	0	0.000 ug/L



Unidentified Peaks

None

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Revision Log

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Sample ID:	6507-2	Operator:	Jeremy Armstrong
Instrument ID:	Varian GC/MS #1	Last Calibration:	4/9/2012 8:17 AM
Acquisition Date:	4/24/2012 4:16 PM	Data File:	...il\042312\6507-2.sms
Calculation Date:	4/24/2012 7:18 PM	Method:	..._cain_040612_jwa.mth
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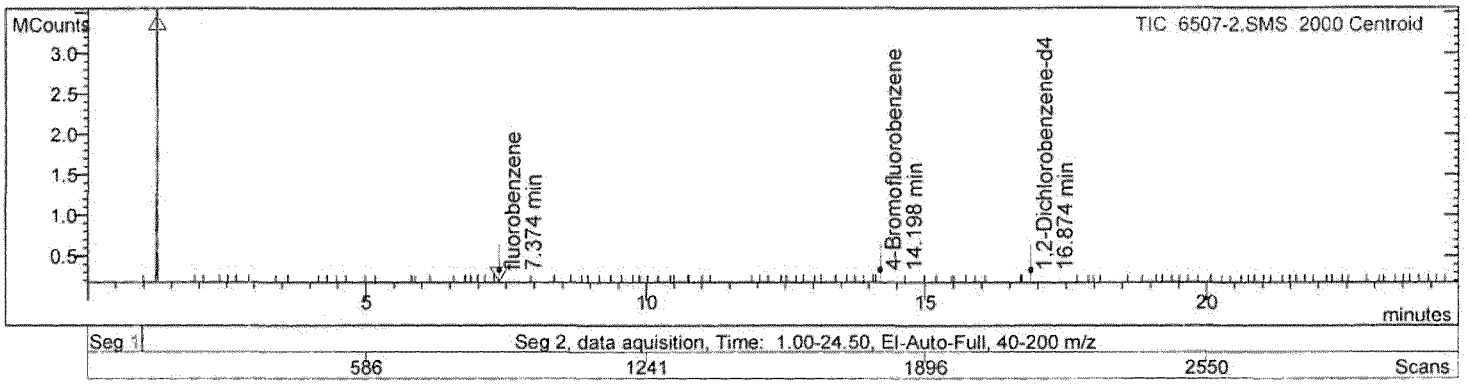
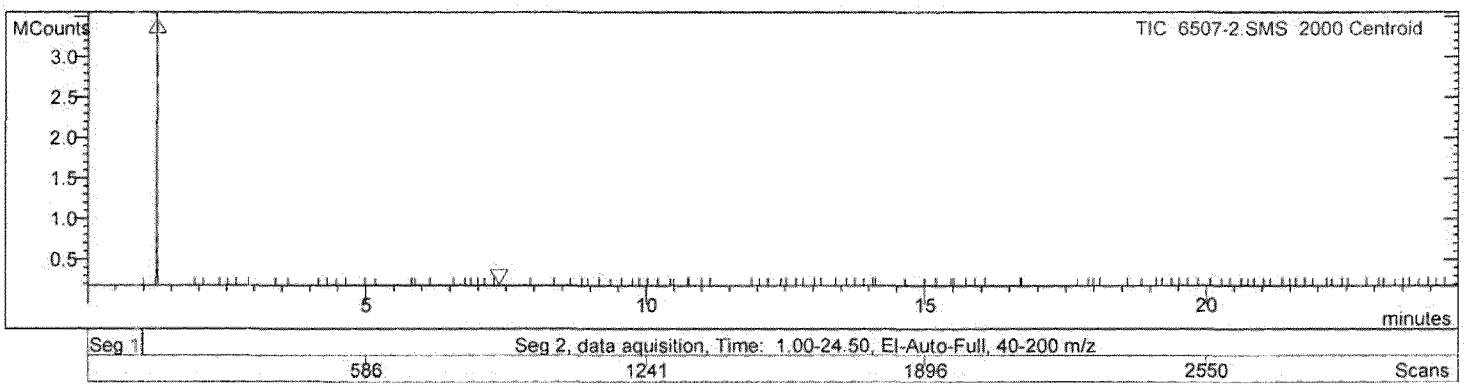
Multiplier:	1.000000		



Target Compounds

#	RT	Peak Name	Res Type	Quan Ions	Area	Amount/RF
1	7.374	fluorobenzene	Id.	96.0	212209	100.000 ug/L
2	2.685	Ethanol	Miss.	45.0+44.0	0	0.000 ug/L
3	3.383	isopropanol	Miss.	45.0+43.0	0	0.000 ug/L
4	5.039	n-propanol	Miss.	59.0+42.0	0	0.000 ug/L
5	6.927	iso-butanol	Miss.	43.1+56.0	0	0.000 ug/L
6	7.939	n-butanol	Miss.	56.0+55.0	0	0.000 ug/L
7	14.198	4-Bromofluorobenzene	Id.	95.0+176.0	138176	104.143 ug/L
8	16.874	1,2-Dichlorobenzene-d4	Id.	115.0	32341	95.752 ug/L



Unidentified Peaks

None

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Revision Log

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Sample ID:	6507-2	Operator:	Jeremy Armstrong
Instrument ID:	Varian GC/MS #1	Last Calibration:	4/16/2012 10:56 AM
Acquisition Date:	4/24/2012 4:16 PM	Data File:	...il\042312\6507-2.sms
Calculation Date:	4/24/2012 7:17 PM	Method:	...312_ccc_cain_jwa.mth
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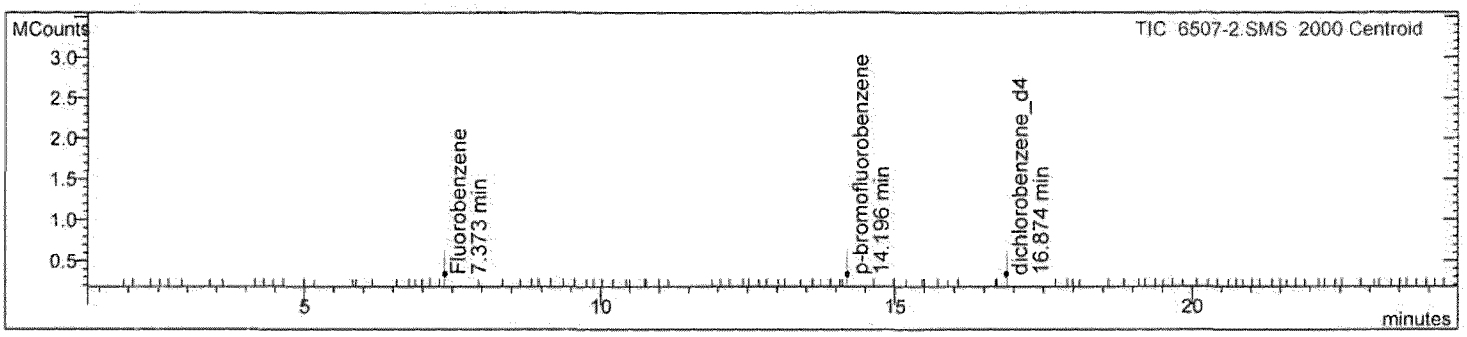
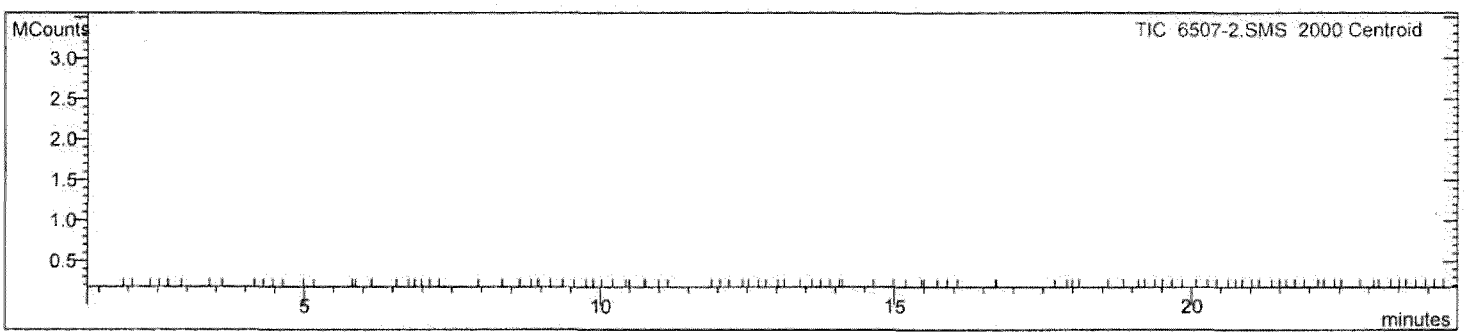
Multiplier:	1.000000		



Target Compounds

#	RT	Peak Name	Res Type	Quan Ions	Area	Amount/RF
1	7.373	Fluorobenzene	Id.	96.0	211937	100.000 ug/L
2	3.160	acetone	Miss.	43.0	0	0.000 ug/L
3	4.000	TBA	Miss.	59.0	0	0.000 ug/L
4	4.194	MTBE	Miss.	73.0	0	0.000 ug/L
5	14.196	p-bromofluorobenzene	Id.	174.0	48289	105.022 ug/L
6	16.874	dichlorobenzene_d4	Id.	115.0	32028	105.129 ug/L



Unidentified Peaks

None

Sample ID:	6507-3	Operator:	Jeremy Armstrong
Instrument ID:	Varian GC/MS #1	Last Calibration:	4/17/2012 10:51 AM
Acquisition Date:	4/24/2012 5:07 PM	Data File:	...il\042312\6507-3.sms
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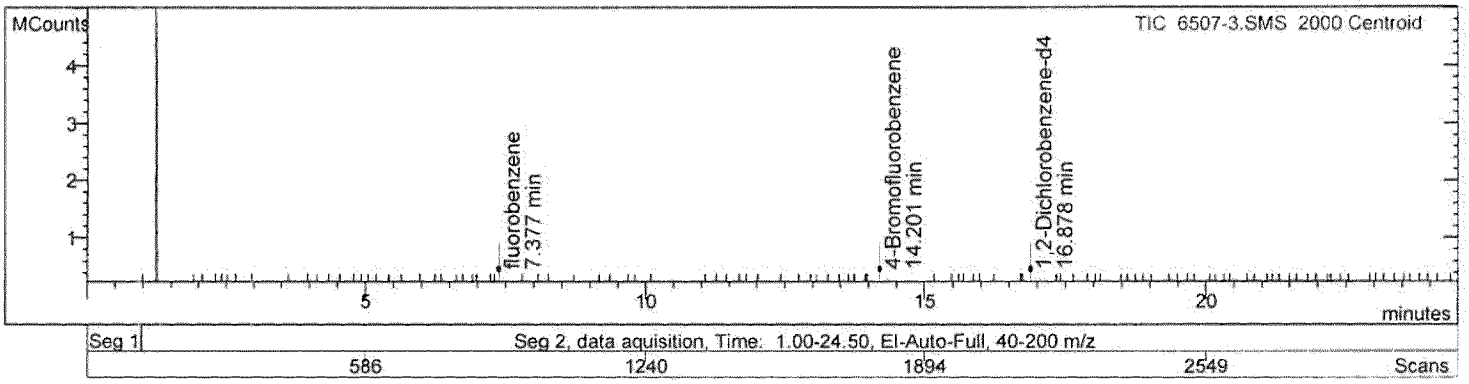
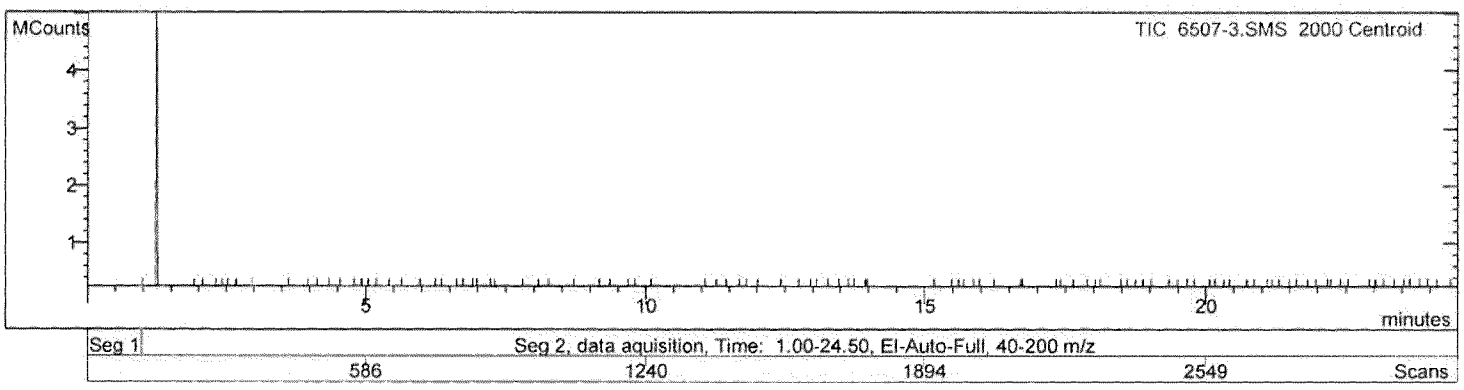


Target Compounds

#	RT	Peak Name	Res Type	Quan Ions	Area	Amount/RF
1	7.377	fluorobenzene	Id.	96.0	225251	100.000 ug/L
2	3.971	TBA	Miss.	59.0+57.0	0	0.000 ug/L
3	4.166	MTBE	Miss.	73.0+57.0	0	0.000 ug/L
4	4.932	di-isopropyl ether (DIPE)	Miss.	45.0+69.0	0	0.000 ug/L
5	5.470	ethyl tert-butyl ether	Miss.	59.0+87.0	0	0.000 ug/L
6	6.907	benzene	Miss.	78.0+77.0	0	0.000 ug/L
7	7.137	tert-amyl methyl ether (T	Miss.	73.0+87.0	0	0.000 ug/L
8	8.057	2,5-dimethylfuran	Miss.	96.0+95.0	0	0.000 ug/L
9	9.908	toluene	Miss.	91.0+92.0	0	0.000 ug/L
10	11.342	1,2-Dibromoethane	Miss.	108.9	0	0.000 ug/L
11	12.413	ethyl benzene	Miss.	91.0+106.0	0	0.000 ug/L
12	12.622	m+p xylene	Miss.	91.0+106.0	0	0.000 ug/L
13	13.281	o-xylene	Miss.	91.0+106.0	0	0.000 ug/L
14	14.201	4-Bromofluorobenzene	Id.	95.0+176.0	144066	101.169 ug/L
15	14.990	1,3,5-Trimethylbenzene	Miss.	105.0+120.0	0	0.000 ug/L
16	15.634	1,2,4-Trimethylbenzene	Miss.	105.0+120.0	0	0.000 ug/L
17	16.363	1,2,3-Trimethylbenzene	Miss.	105.0+120.0	0	0.000 ug/L
18	16.878	1,2-Dichlorobenzene-d4	Id.	115.0	34187	98.478 ug/L
19	19.901	Naphthalene	Miss.	128.0+127.0	0	0.000 ug/L



Unidentified Peaks

None

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Revision Log

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Sample ID:	6507-3	Operator:	Jeremy Armstrong
Instrument ID:	Varian GC/MS #1	Last Calibration:	4/9/2012 8:17 AM
Acquisition Date:	4/24/2012 5:07 PM	Data File:	...il\042312\6507-3.sms
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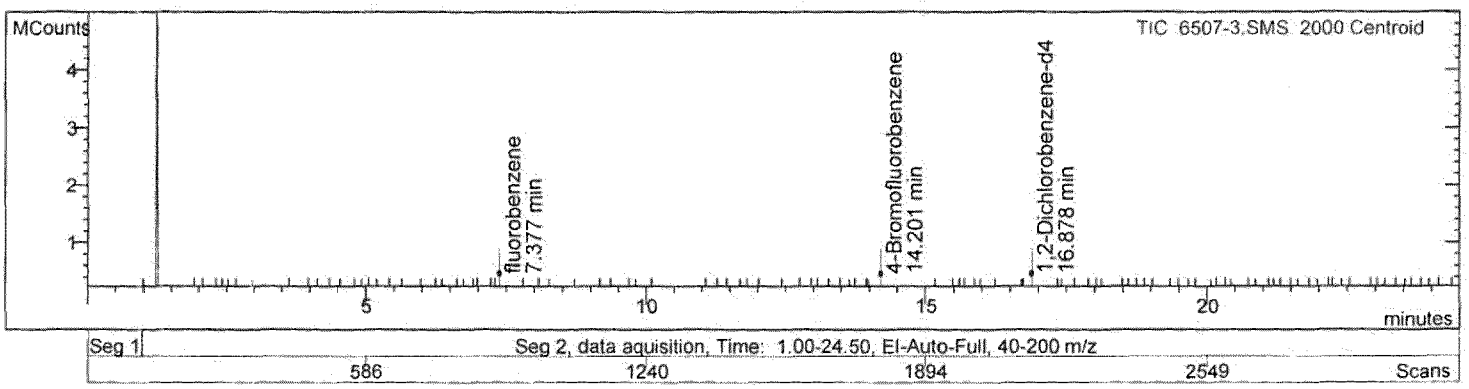
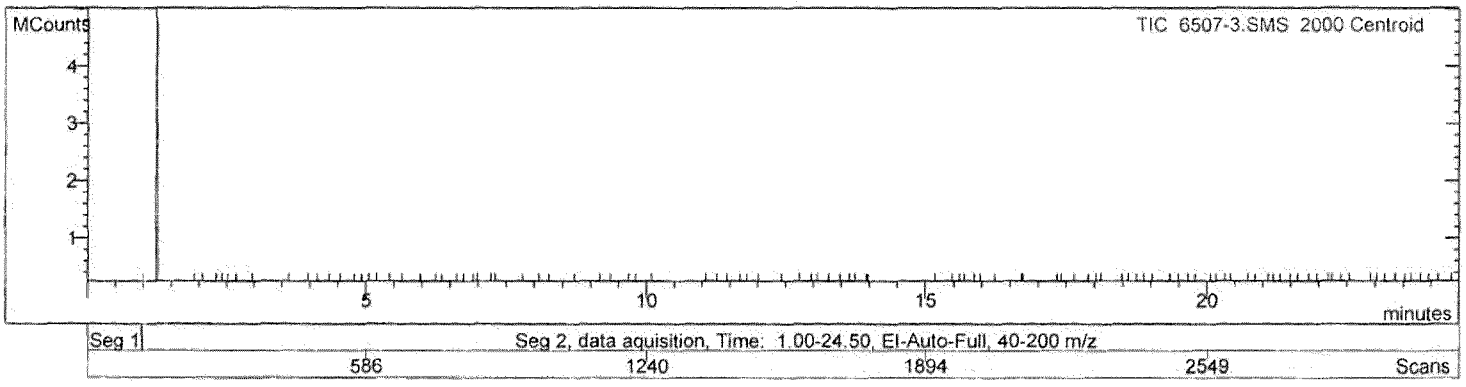


Target Compounds

#	RT	Peak Name	Res Type	Quan Ions	Area	Amount/RF
1	7.377	fluorobenzene	Id.	96.0	225384	100.000 ug/L
2	2.685	Ethanol	Miss.	45.0+44.0	0	0.000 ug/L
3	3.383	isopropanol	Miss.	45.0+43.0	0	0.000 ug/L
4	5.039	n-propanol	Miss.	59.0+42.0	0	0.000 ug/L
5	6.927	iso-butanol	Miss.	43.1+56.0	0	0.000 ug/L
6	7.939	n-butanol	Miss.	56.0+55.0	0	0.000 ug/L
7	14.201	4-Bromofluorobenzene	Id.	95.0+176.0	144198	102.329 ug/L
8	16.878	1,2-Dichlorobenzene-d4	Id.	115.0	34187	95.0298 ug/L



Unidentified Peaks

None

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Revision Log

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Sample ID:	6507-3	Operator:	Jeremy Armstrong
Instrument ID:	Varian GC/MS #1	Last Calibration:	4/16/2012 10:56 AM
Acquisition Date:	4/24/2012 5:07 PM	Data File:	...il\042312\6507-3.sms
Calculation Date:	4/24/2012 7:17 PM	Method:	...312_ccc_cain_jwa.mth
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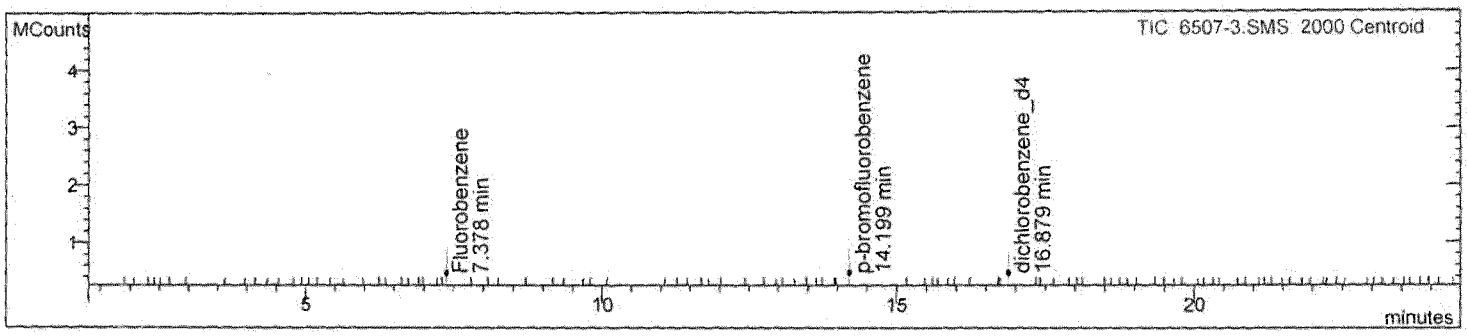
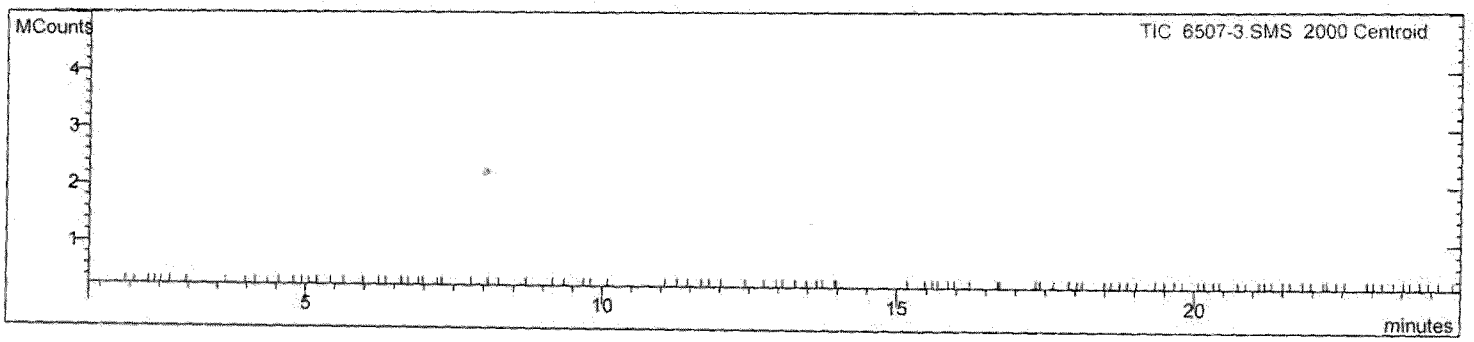


Target Compounds

#	RT	Peak Name	Res Type	Quan Ions	Area	Amount/RF
1	7.378	Fluorobenzene	Id.	96.0	224923	100.000 ug/L
2	3.160	acetone	Miss.	43.0	0	0.000 ug/L
3	4.000	TBA	Miss.	59.0	0	0.000 ug/L
4	4.194	MTBE	Miss.	73.0	0	0.000 ug/L
5	14.199	p-bromofluorobenzene	Id.	174.0	50615	103.724 ug/L
6	16.879	dichlorobenzene_d4	Id.	115.0	33846	104.681 ug/L



Unidentified Peaks

None

Sample ID:	6507-4	Operator:	Jeremy Armstrong
Instrument ID:	Varian GC/MS #1	Last Calibration:	4/17/2012 10:51 AM
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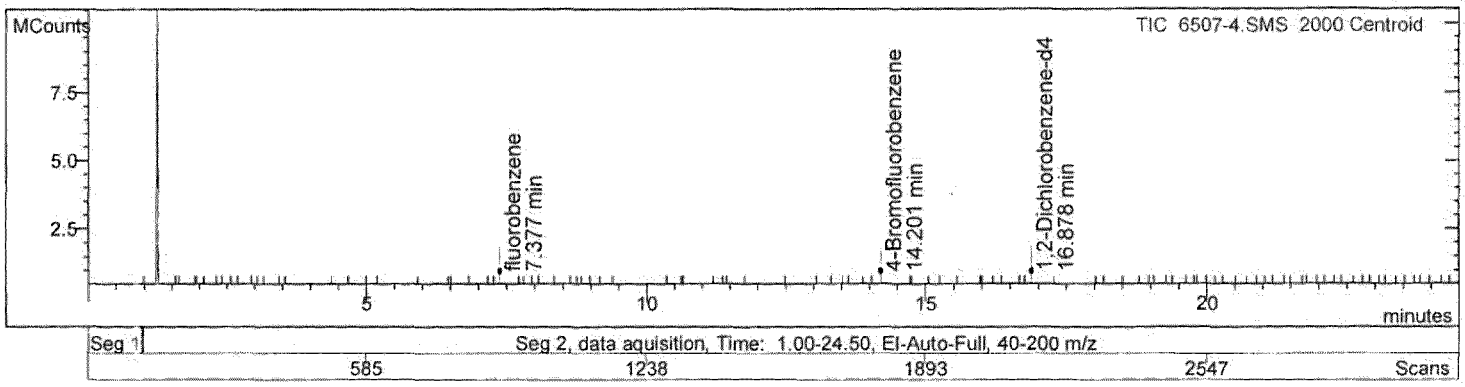
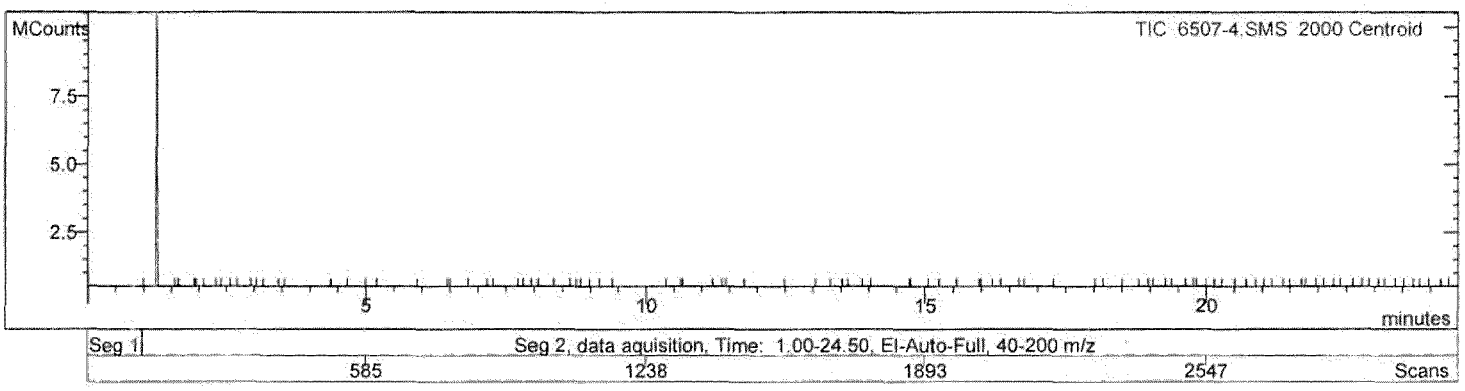


Target Compounds

#	RT	Peak Name	Res Type	Quan Ions	Area	Amount/RF
1	7.377	fluorobenzene	Id.	96.0	185800	100.000 ug/L
2	3.971	TBA	Miss.	59.0+57.0	0	0.000 ug/L
3	4.166	MTBE	Miss.	73.0+57.0	0	0.000 ug/L
4	4.932	di-isopropyl ether (DIPE)	Miss.	45.0+69.0	0	0.000 ug/L
5	5.470	ethyl tert-butyl ether	Miss.	59.0+87.0	0	0.000 ug/L
6	6.907	benzene	Miss.	78.0+77.0	0	0.000 ug/L
7	7.137	tert-amyl methyl ether (T	Miss.	73.0+87.0	0	0.000 ug/L
8	8.057	2,5-dimethylfuran	Miss.	96.0+95.0	0	0.000 ug/L
9	9.908	toluene	Miss.	91.0+92.0	0	0.000 ug/L
10	11.342	1,2-Dibromoethane	Miss.	108.9	0	0.000 ug/L
11	12.413	ethyl benzene	Miss.	91.0+106.0	0	0.000 ug/L
12	12.622	m+p xylene	Miss.	91.0+106.0	0	0.000 ug/L
13	13.281	o-xylene	Miss.	91.0+106.0	0	0.000 ug/L
14	14.201	4-Bromofluorobenzene	Id.	95.0+176.0	106616	90.767 ug/L
15	14.990	1,3,5-Trimethylbenzene	Miss.	105.0+120.0	0	0.000 ug/L
16	15.634	1,2,4-Trimethylbenzene	Miss.	105.0+120.0	0	0.000 ug/L
17	16.363	1,2,3-Trimethylbenzene	Miss.	105.0+120.0	0	0.000 ug/L
18	16.878	1,2-Dichlorobenzene-d4	Id.	115.0	23846	83.275 ug/L
19	19.901	Naphthalene	Miss.	128.0+127.0	0	0.000 ug/L



Unidentified Peaks

None

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Revision Log

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Sample ID:	6507-4	Operator:	Jeremy Armstrong
Instrument ID:	Varian GC/MS #1	Last Calibration:	4/9/2012 8:17 AM
Acquisition Date:	4/24/2012 5:58 PM	Data File:	...il\042312\6507-4.sms
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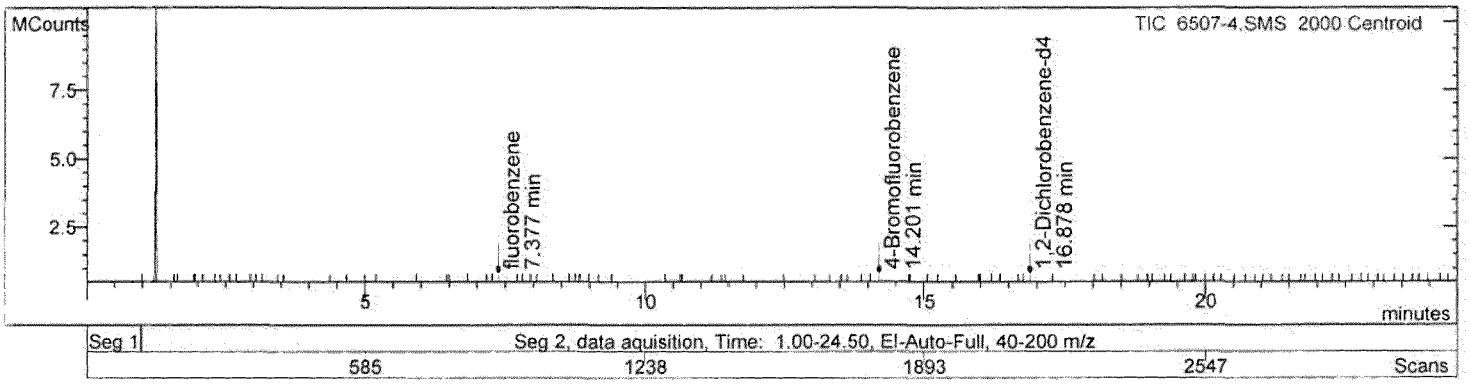
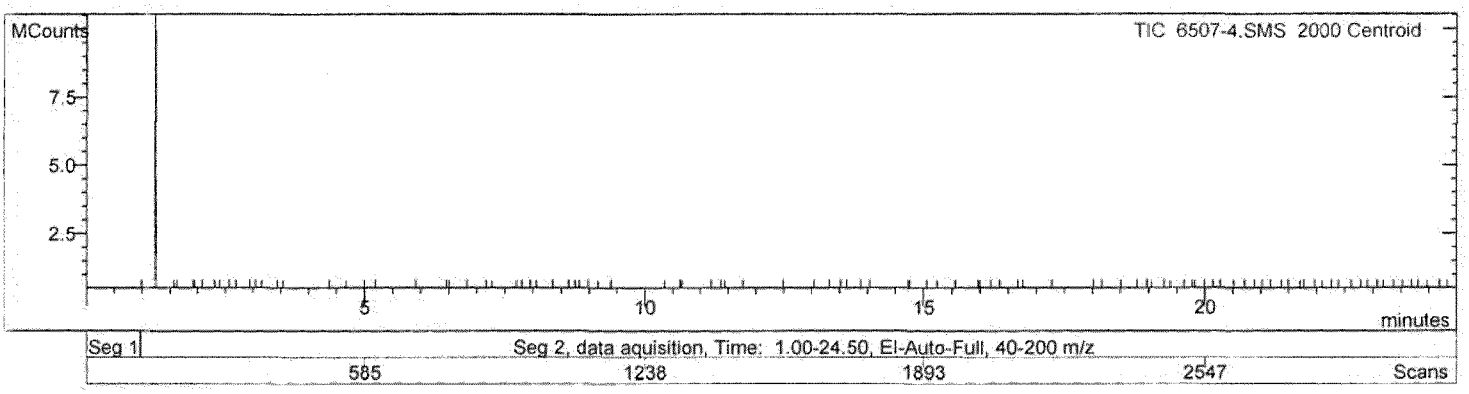


Target Compounds

#	RT	Peak Name	Res Type	Quan Ions	Area	Amount/RF
1	7.377	fluorobenzene	Id.	96.0	185787	100.000 ug/L
2	2.685	Ethanol	Miss.	45.0+44.0	0	0.000 ug/L
3	3.383	isopropanol	Miss.	45.0+43.0	0	0.000 ug/L
4	5.039	n-propanol	Miss.	59.0+42.0	0	0.000 ug/L
5	6.927	iso-butanol	Miss.	43.1+56.0	0	0.000 ug/L
6	7.939	n-butanol	Miss.	56.0+55.0	0	0.000 ug/L
7	14.201	4-Bromofluorobenzene	Id.	95.0+176.0	106616	91.784 ug/L
8	16.878	1,2-Dichlorobenzene-d4	Id.	115.0	23846	80.640 ug/L



Unidentified Peaks

None

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Revision Log

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Sample ID:	6507-4	Operator:	Jeremy Armstrong
Instrument ID:	Varian GC/MS #1	Last Calibration:	4/16/2012 10:56 AM
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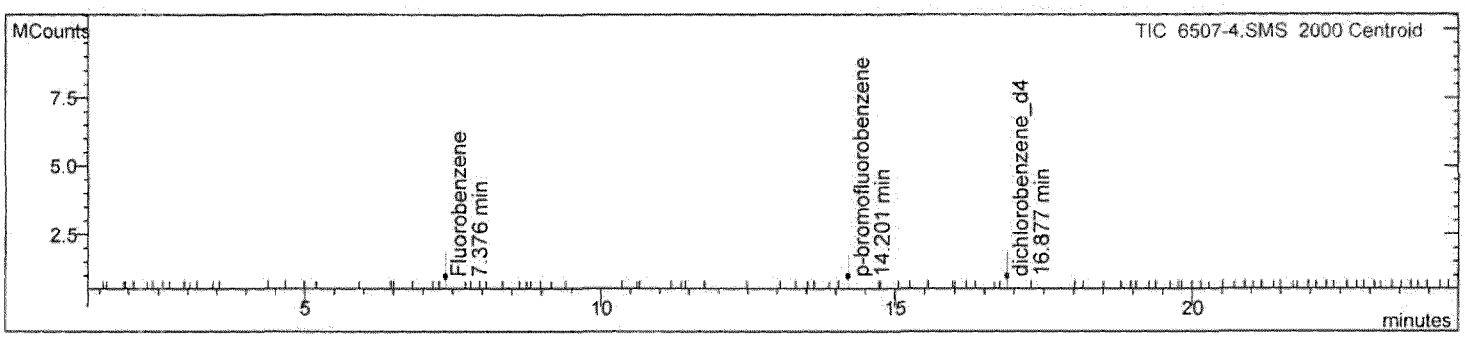
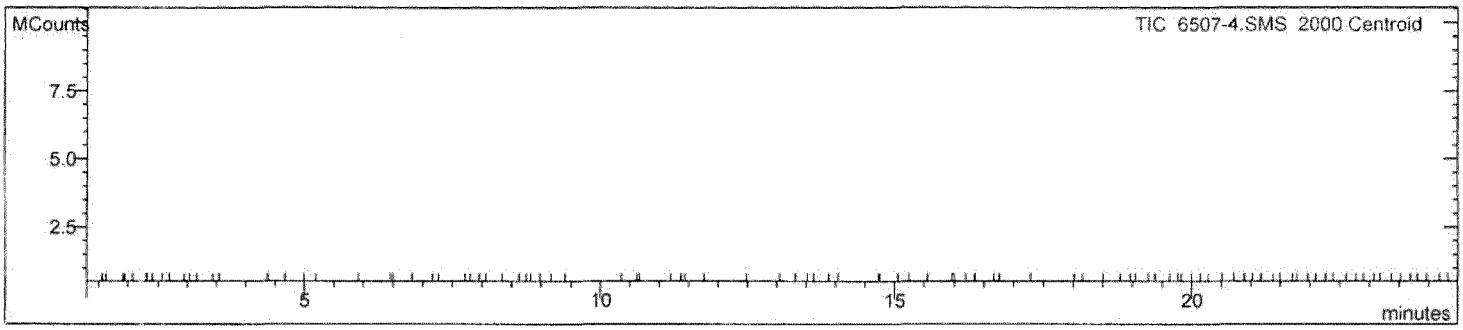


Target Compounds

#	RT	Peak Name	Res Type	Quan Ions	Area	Amount/RF
1	7.376	Fluorobenzene	Id.	96.0	185450	100.000 ug/L
2	3.160	acetone	Miss.	43.0	0	0.000 ug/L
3	4.000	TBA	Miss.	59.0	0	0.000 ug/L
4	4.194	MTBE	Miss.	73.0	0	0.000 ug/L
5	14.201	p-bromofluorobenzene	Id.	174.0	36989	91.935 ug/L
6	16.877	dichlorobenzene_d4	Id.	115.0	23607	88.555 ug/L



Unidentified Peaks

None

Sample ID:	6507-4	Operator:	Jeremy Armstrong
Instrument ID:	Varian GC/MS #1	Last Calibration:	4/20/2012 12:37 PM
Acquisition Date:	4/24/2012 5:58 PM	Data File:	...il\042312\6507-4.sms
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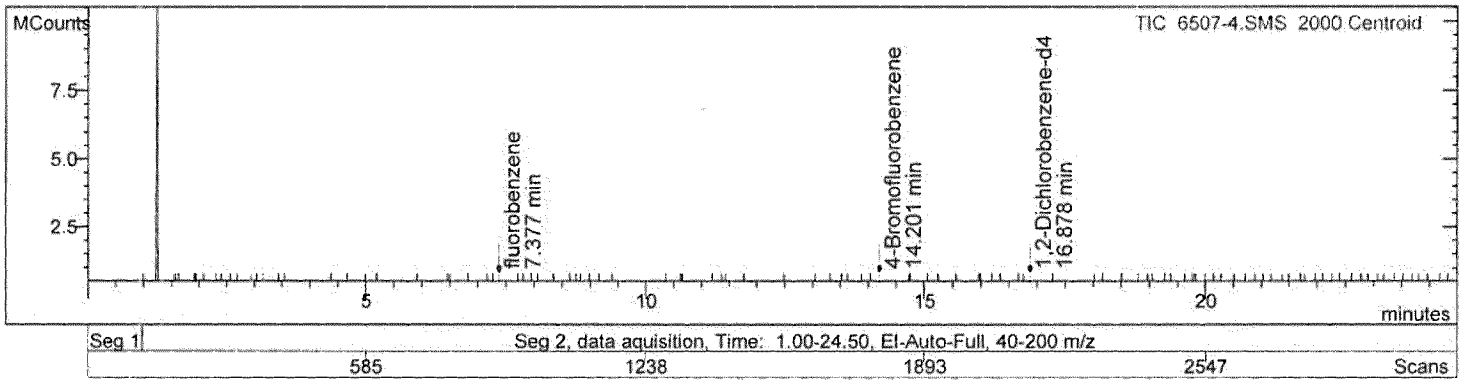
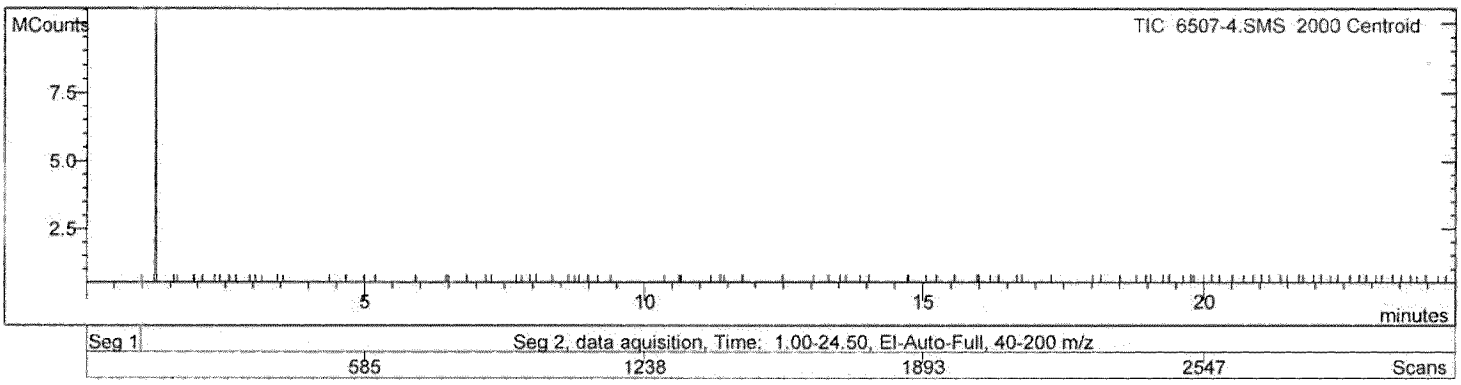


Target Compounds

#	RT	Peak Name	Res Type	Quan Ions	Area	Amount/RF
1	7.377	fluorobenzene	Id.	96.0	185787	100.000 ug/L
2	1.677	Vinyl Chloride	Miss.	62.0+64.0	0	0.000 ug/L
3	3.075	1,1-Dichloroethene	Miss.	61.0+96.0	0	0.000 ug/L
4	3.770	Methylene Chloride	Miss.	49.0+84.0	0	0.000 ug/L
5	4.176	trans-1,2-dichloroethene	Miss.	61.0+96.0	0	0.000 ug/L
6	4.793	1,1-dichloroethane	Miss.	63.0+65.0	0	0.000 ug/L
7	5.661	cis-1,2-dichloroethene	Miss.	61.0+96.0	0	0.000 ug/L
8	6.156	chloroform	Miss.	83.0+85.0	0	0.000 ug/L
9	6.391	1,1,1-trichloroethane	Miss.	97.0+61.0	0	0.000 ug/L
10	6.634	carbon tetrachloride	Miss.	117.0+119.0	0	0.000 ug/L
11	6.980	1,2-dichloroethane	Miss.	62.0+64.0	0	0.000 ug/L
12	7.939	trichloroethene	Miss.	95.0+132.0	0	0.000 ug/L
13	10.613	1,1,2-Trichloroethane	Miss.	96.8	0	0.000 ug/L
14	10.817	tetrachloroethene	Miss.	166.0+129.0	0	0.000 ug/L
15	12.225	chlorobenzene	Miss.	112.0+77.0	0	0.000 ug/L
16	14.201	4-Bromofluorobenzene	Id.	95.0+176.0	106616	92.676 ug/L
17	16.106	1,3-dichlorobenzene	Miss.	146.1+111.0	0	0.000 ug/L
18	16.271	1,4-dichlorobenzene	Miss.	146.1+111.0	0	0.000 ug/L
19	16.878	1,2-Dichlorobenzene-d4	Id.	115.0	23846	90.648 ug/L
20	16.916	1,2-dichlorobenzene	Miss.	146.1	0	0.000 ug/L



Unidentified Peaks

None

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Revision Log

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Print Date: 24 Apr 2012 19:13:45

Chromatogram Plot

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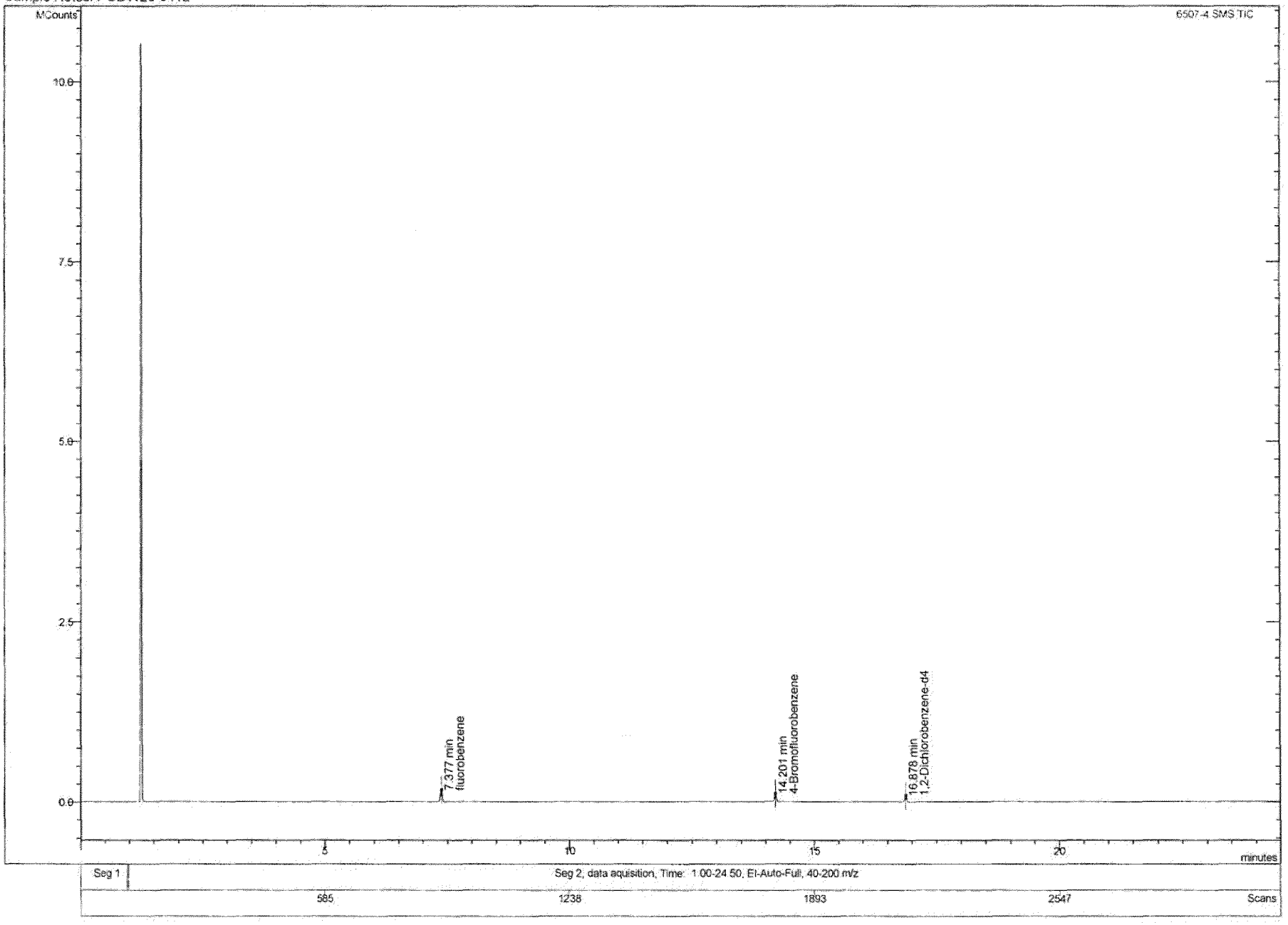
Sample: 6507-4
Scan Range: 1 - 3127 Time Range: 0.00 - 24.49 min.
Sample Notes: PGDW20-0412

Operator: Jeremy Armstrong
Date: 4/24/2012 5:58 PM



EPAPAV0043736