LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM032819\
Data File : BM019567.D
Acq On : 29 Mar 2019 07:04
Operator : JU/SJ
Sample : K2063-16DL 10X
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : DB6S4DL

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM032219MA.M
Title : SVOA CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.375	63	66	74	rVB	132530	181076	0.34%	0.088%
2	3.628	101	109	121	rVB	119930	245630	0.46%	0.119%
3	3.828	136	143	164	rBV2	32661018	52192541	98.31%	25.320%
4	3.969	164	167	174	rBV2	63202	97014	0.18%	0.047%
5	4.028	174	177	179	rBV	46501	61067	0.12%	0.030%
6	4.116	187	192	210	rVB3	54526	127205	0.24%	0.062%
7	4.599	269	274	286	rBV	71126	144162	0.27%	0.070%
8	4.705	288	292	298	rVV2	45456	76348	0.14%	0.037%
9	4.769	298	303	308	rVV	127990	194973	0.37%	0.095%
10	4.822	308	312	322	rVB	100129	176165	0.33%	0.085%
11	5.110	354	361	372	rBV	9828111	14581423	27.47%	7.074%
12	5.246	376	384	401	rVB	24872833	53088157	100.00%	25.755%
13	5.604	436	445	455	rVV	17368109	24924275	46.95%	12.091%
14	5.687	455	459	462	rVV	44791	71956	0.14%	0.035%
15	5.722	462	465	472	rVB	36469	55305	0.10%	0.027%
16	6.069	515	524	537	rBV	355442	612490	1.15%	0.297%
17	6.251	550	555	564	rVB2	47440	88144	0.17%	0.043%
18	6.551	600	606	615	rVB	978305	1407395	2.65%	0.683%
19	6.663	615	625	630	rBV	4144835	6048850	11.39%	2.934%
20	6.716	630	634	641	rVV	1649258	2430564	4.58%	1.179%
21	6.793	641	647	655	rVB	1712616	2506949	4.72%	1.216%
22	6.957	665	675	686	rBV	1770307	2675321	5.04%	1.298%
23	7.116	697	702	708	rBV	51306	89258	0.17%	0.043%
24	7.216	712	719	729	rVB	7034111	10133075	19.09%	4.916%
25	7.575	774	780	790	rBV2	594443	1067464	2.01%	0.518%
26	7.675	790	797	806	rVB	1813414	2752831	5.19%	1.335%
27	7.934	835	841	848	rVB	1447479	2290428	4.31%	1.111%
28	8.045	855	860	865	rBV	129788	205762	0.39%	0.100%
29	8.104	865	870	879	rVB	344040	537667	1.01%	0.261%
30	8.198	879	886	891	rBV	445740	704470	1.33%	0.342%
31	8.357	908	913	921	rVB	94585	149368	0.28%	0.072%
32	8.510	930	939	943	rBV	243120	409808	0.77%	0.199%
33	8.557	943	947	952	rVV	205535	357047	0.67%	0.173%
34	8.616	952	957	960	rVV	136448	242908	0.46%	0.118%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM032819\
Data File : BM019567.D
Acq On : 29 Mar 2019 07:04
Operator : JU/SJ
Sample : K2063-16DL 10X
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : DB6S4DL

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM032219MA.M
Title : SVOA CALIBRATION

35	8.657	960	964	972	rVV2	406121	659709	1.24%	0.320%
36	8.740	972	978	991	rVB3	220905	418199	0.79%	0.203%
37	8.992	1015	1021	1028	rBV	108099	183276	0.35%	0.089%
38	9.181	1042	1053	1058	rBV	212608	332086	0.63%	0.161%
39	9.240	1058	1063	1071	rVB	316957	464460	0.87%	0.225%
40	9.469	1097	1102	1111	rVB3	51192	121623	0.23%	0.059%
41	9.598	1119	1124	1133	rVB	240910	402426	0.76%	0.195%
42	9.745	1136	1149	1158	rBV2	484011	896781	1.69%	0.435%
43	9.987	1183	1190	1197	rBV4	78434	198278	0.37%	0.096%
44	10.081	1201	1206	1214	rVB	77963	143722	0.27%	0.070%
45	10.316	1240	1246	1247	rBV	84592	146290	0.28%	0.071%
46	10.357	1247	1253	1257	rVV	848316	1483619	2.79%	0.720%
47	10.410	1257	1262	1268	rVB	1923651	3088381	5.82%	1.498%
48	10.534	1276	1283	1295	rVB2	46226	109777	0.21%	0.053%
49	10.881	1336	1342	1353	rBV2	59264	116039	0.22%	0.056%
50	11.175	1385	1392	1400	rBV	115098	240804	0.45%	0.117%
51	11.610	1460	1466	1473	rVB	39564	66864	0.13%	0.032%
52	11.786	1492	1496	1505	rVB2	40403	80006	0.15%	0.039%
53	12.033	1530	1538	1548	rBV	400876	673627	1.27%	0.327%
54	12.210	1563	1568	1571	rBV2	83741	134390	0.25%	0.065%
55	12.251	1571	1575	1582	rVB	206007	336234	0.63%	0.163%
56	12.404	1591	1601	1615	rBV	194322	400280	0.75%	0.194%
57	12.698	1646	1651	1660	rVB3	47084	81867	0.15%	0.040%
58	13.222	1735	1740	1745	rBV	51285	82219	0.15%	0.040%
59	13.422	1770	1774	1780	rVB	64339	98136	0.18%	0.048%
60	13.663	1810	1815	1822	rVB	139016	208398	0.39%	0.101%
61	13.927	1855	1860	1873	rVB	160797	252085	0.47%	0.122%
62	14.239	1906	1913	1921	rBV2	1314288	1949655	3.67%	0.946%
63	15.239	2077	2083	2088	rBV	207620	305073	0.57%	0.148%
64	16.998	2376	2382	2394	rBV	1661905	2314520	4.36%	1.123%
65	17.098	2394	2399	2410	rVB2	201516	325385	0.61%	0.158%
66	19.404	2786	2791	2805	rBV	291119	441820	0.83%	0.214%
67	21.203	3092	3097	3108	rBV	2544052	3305319	6.23%	1.604%
68	22.221	3267	3270	3275	rVB3	64772	75212	0.14%	0.036%
69	22.709	3350	3353	3358	rVB2	106324	147123	0.28%	0.071%
70	23.268	3442	3448	3455	rVB3	394901	730432	1.38%	0.354%
71	23.409	3465	3472	3484	rVB2	2007343	3780707	7.12%	1.834%

LSC Area Percent Report

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Data File : BM019567.D
Acq On : 29 Mar 2019 07:04
Operator : JU/SJ
Sample : K2063-16DL 10X
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : DB6S4DL

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM032219MA.M
Title : SVOA CALIBRATION

72	23.886	3549	3553	3560	rVB2	105882	197614	0.37%	0.096%
73	24.609	3672	3676	3685	rVB3	100653	211153	0.40%	0.102%

Sum of corrected areas: 206130685

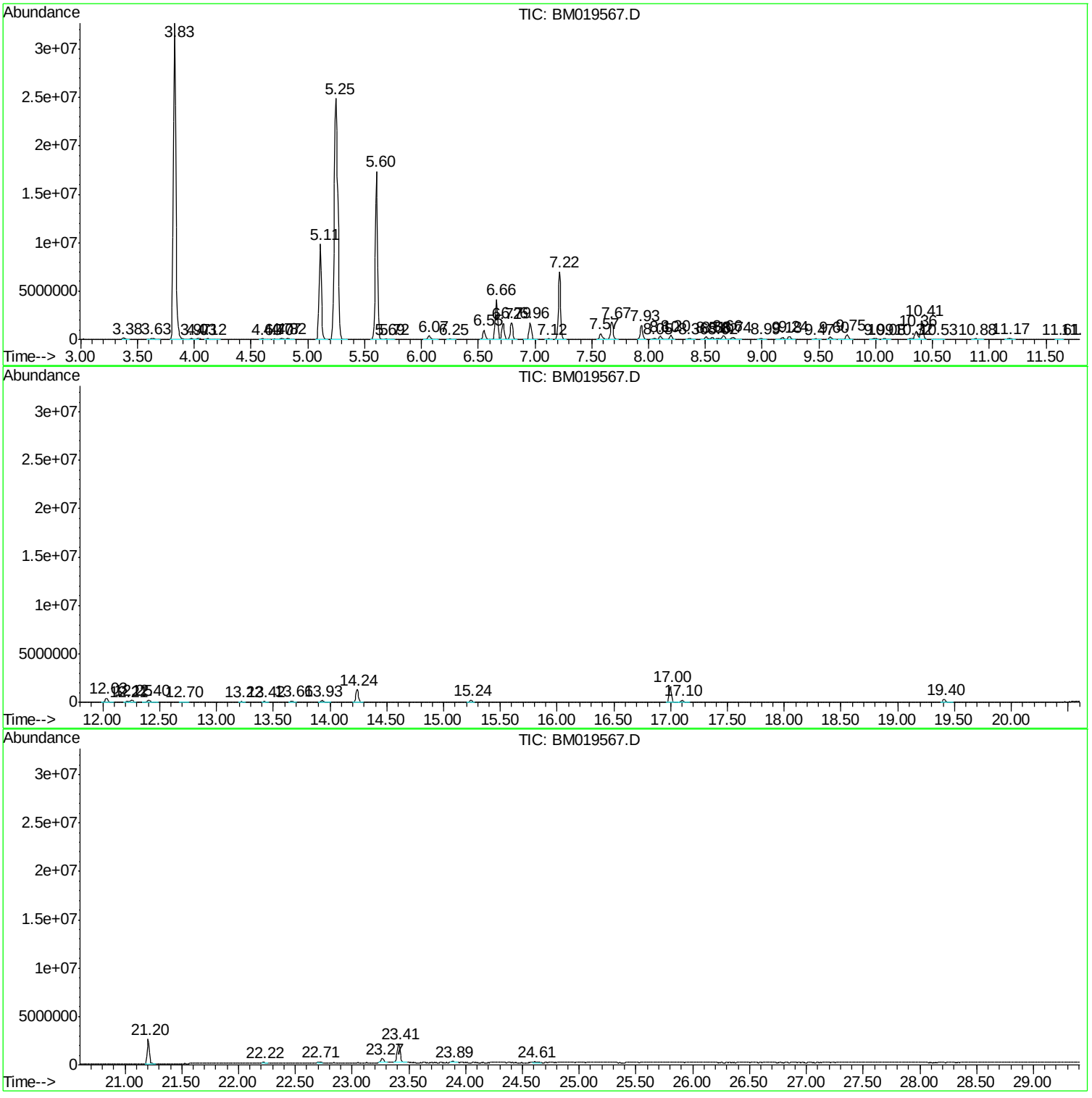
LSC Report - Integrated Chromatogram

Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM032819\
Data File : BM019567.D
Acq On : 29 Mar 2019 07:04
Operator : JU/SJ
Sample : K2063-16DL 10X
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : DB6S4DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM032219MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



Library Search Compound Report

Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM032819\
Data File : BM019567.D
Acq On : 29 Mar 2019 07:04
Operator : JU/SJ
Sample : K2063-16DL 10X
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : DB6S4DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM032219MA.M
Quant Title : SVOA CALIBRATION

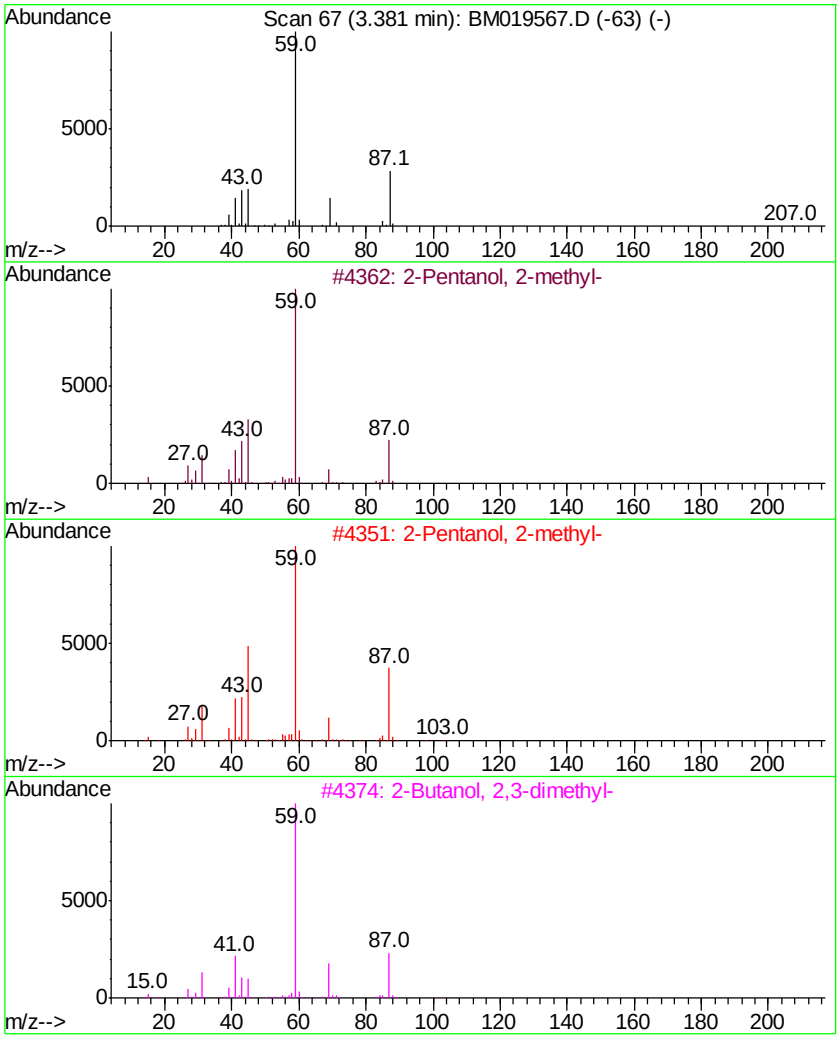
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

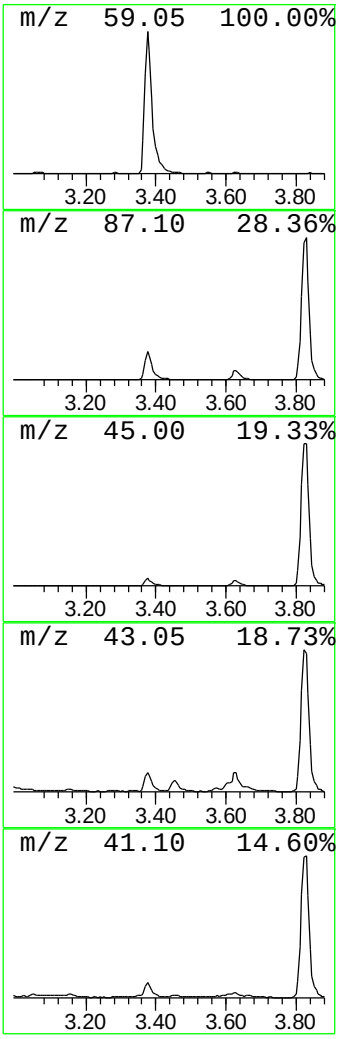
****

Peak Number 1 2-Pentanol, 2-methyl- Concentration Rank 27

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.38	3.39 ng/ul	181076	1,4-Dichlorobenzene-d4	7.57

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanol, 2-methyl-	102	C6H14O	000590-36-3	83
2	2-Pentanol, 2-methyl-	102	C6H14O	000590-36-3	74
3	2-Butanol, 2,3-dimethyl-	102	C6H14O	000594-60-5	72
4	2-Butanol, 2,3-dimethyl-	102	C6H14O	000594-60-5	64
5	3-Pentanol, 2,2-dimethyl-	116	C7H16O	003970-62-5	64





Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM032819\
Data File : BM019567.D
Acq On : 29 Mar 2019 07:04
Operator : JU/SJ
Sample : K2063-16DL 10X
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : DB6S4DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM032219MA.M
Quant Title : SVOA CALIBRATION

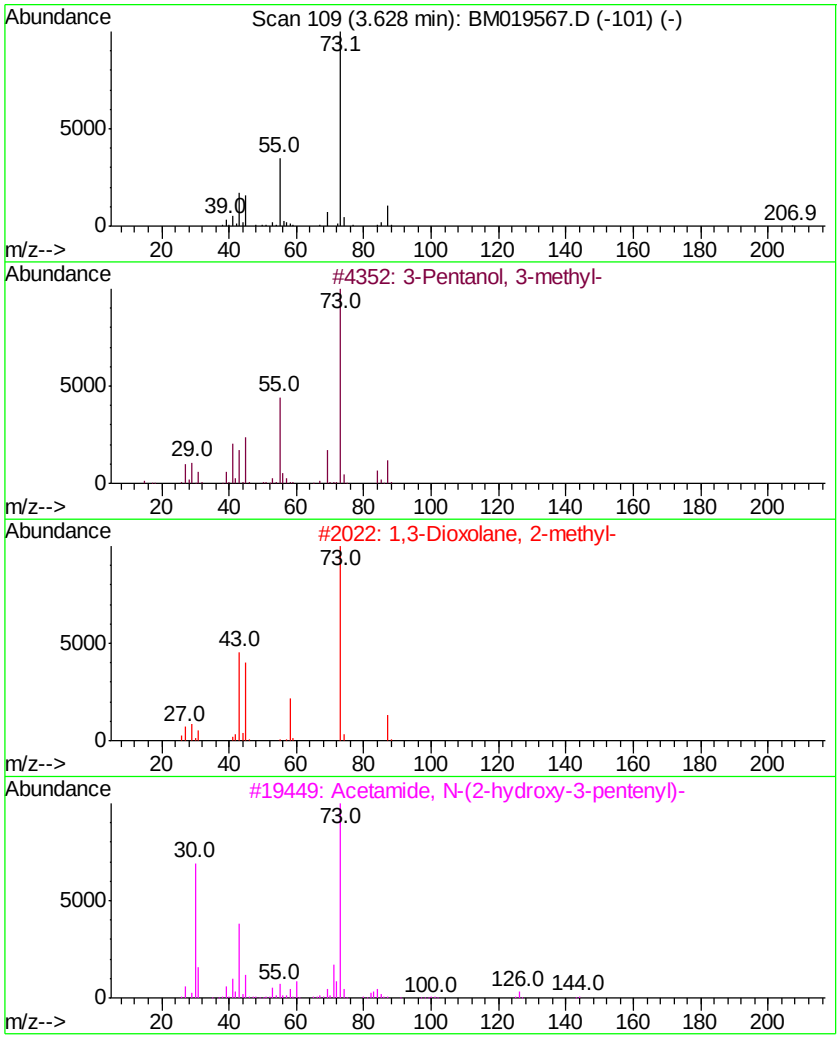
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

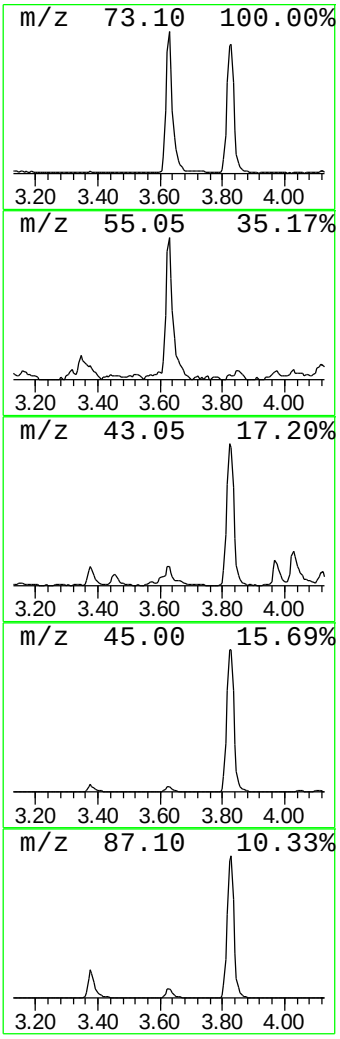
****

Peak Number 2 3-Pentanol, 3-methyl- Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.63	4.60 ng/ul	245630	1,4-Dichlorobenzene-d4	7.57

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Pentanol, 3-methyl-	102	C6H14O	000077-74-7	50
2	1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	42
3	Acetamide, N-(2-hydroxy-3-penten...	143	C7H13NO2	093393-95-4	36
4	3-Pentanol, 3-methyl-	102	C6H14O	000077-74-7	33
5	Octanal, 7-methoxy-3,7-dimethyl-	186	C11H22O2	003613-30-7	28





Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM032819\
Data File : BM019567.D
Acq On : 29 Mar 2019 07:04
Operator : JU/SJ
Sample : K2063-16DL 10X
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : DB6S4DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM032219MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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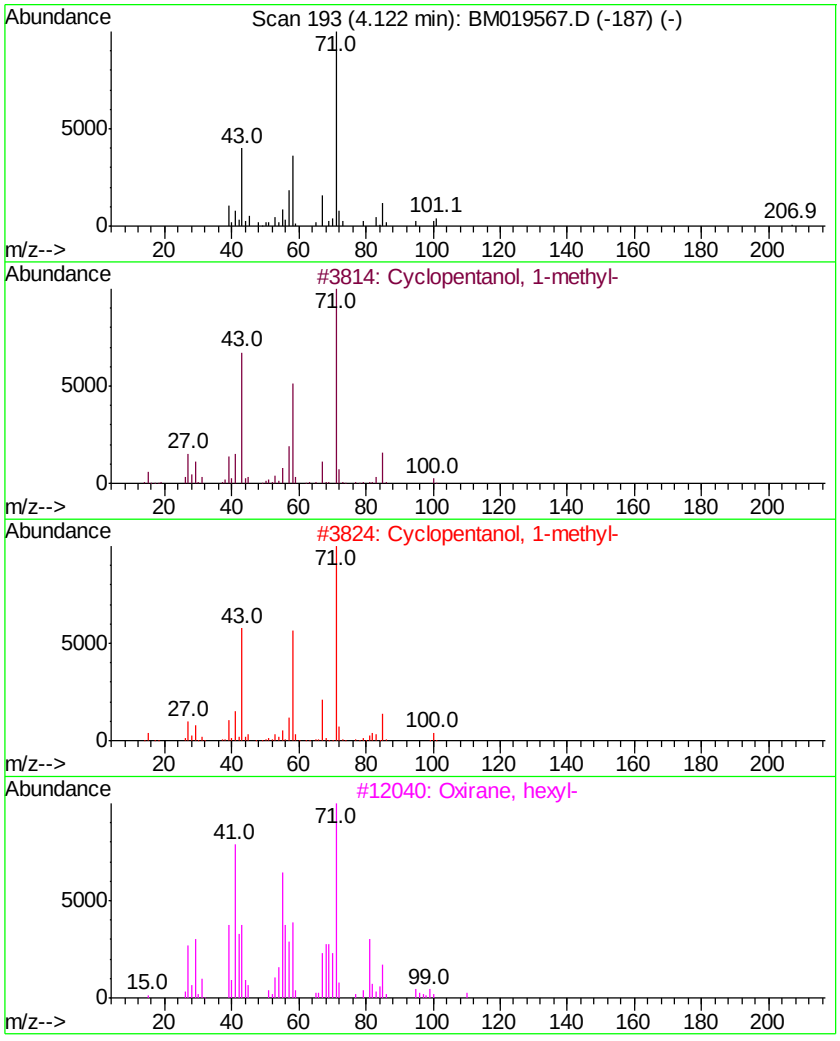
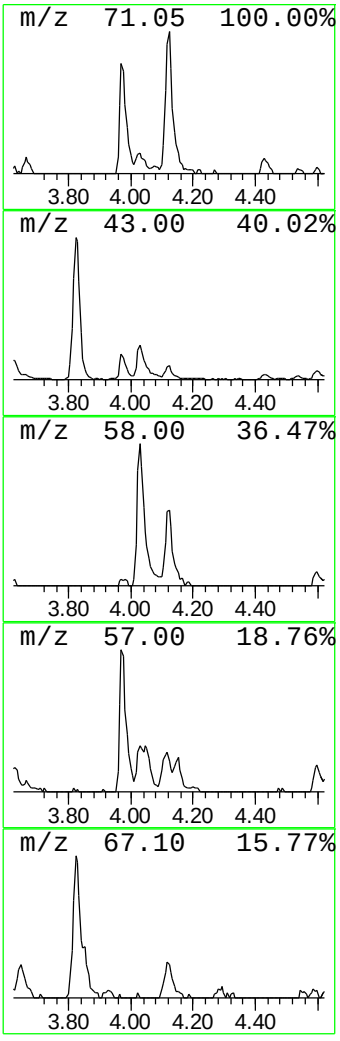
Peak Number 4 Cyclopentanol, 1-methyl- Concentration Rank 33

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.12	2.38 ng/ul	127205	1,4-Dichlorobenzene-d4	7.57

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentanol, 1-methyl-	100	C6H12O	001462-03-9	78
2	Cyclopentanol, 1-methyl-	100	C6H12O	001462-03-9	64
3	Oxirane, hexyl-	128	C8H16O	002984-50-1	64
4	Oxirane, butyl-	100	C6H12O	001436-34-6	50
5	Oxirane, pentyl-	114	C7H14O	005063-65-0	43





Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM032819\
Data File : BM019567.D
Acq On : 29 Mar 2019 07:04
Operator : JU/SJ
Sample : K2063-16DL 10X
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : DB6S4DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM032219MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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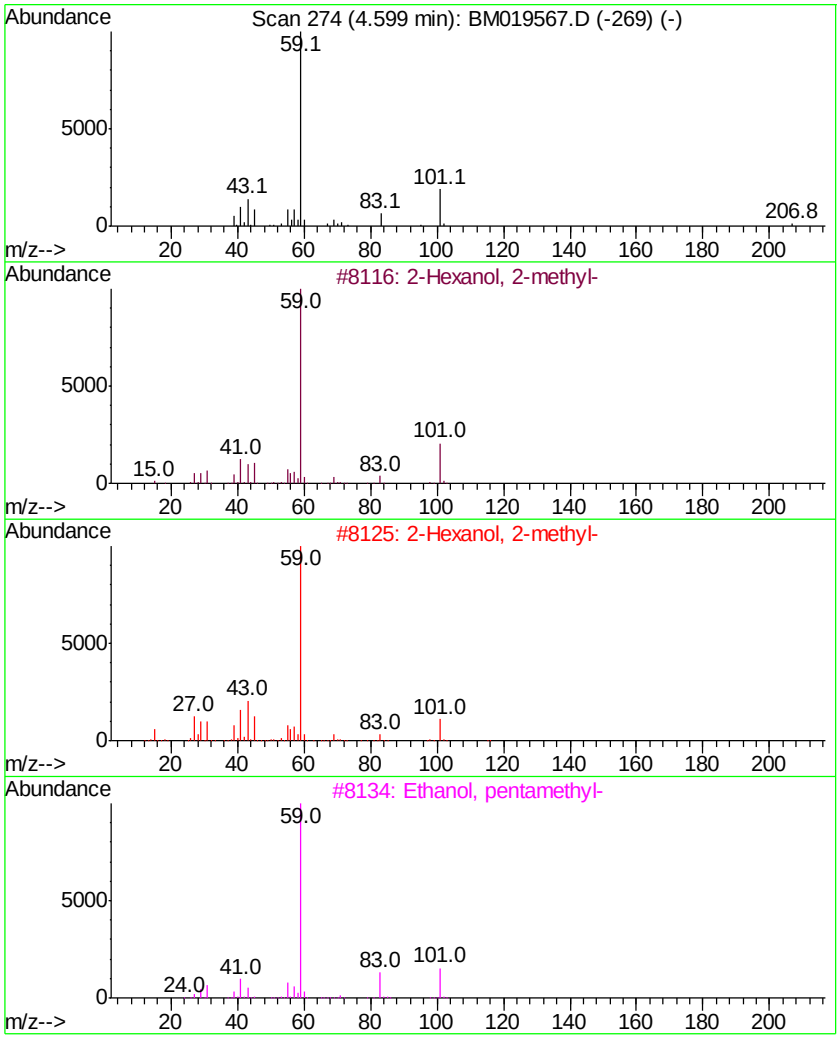
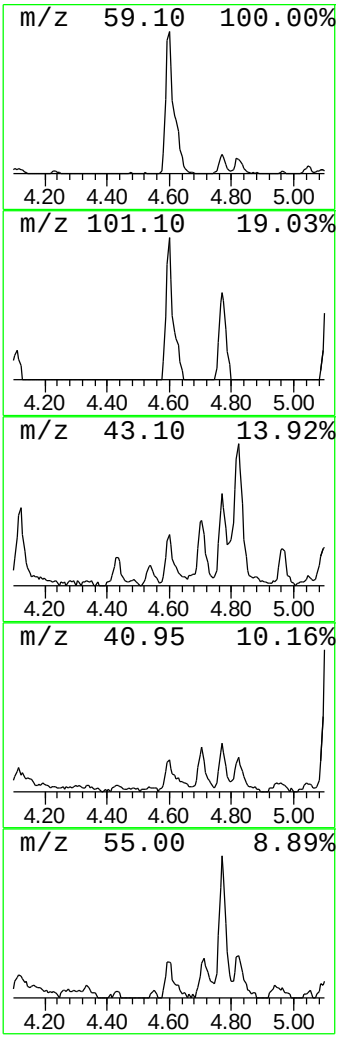
Peak Number 5 2-Hexanol, 2-methyl- Concentration Rank 31

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.60	2.70 ng/ul	144162	1,4-Dichlorobenzene-d4	7.57

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	83
2	2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	78
3	Ethanol, pentamethyl-	116	C7H16O	000594-83-2	78
4	2-Pentanol, 2,3-dimethyl-	116	C7H16O	004911-70-0	74
5	2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	64





Library Search Compound Report

Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM032819\
Data File : BM019567.D
Acq On : 29 Mar 2019 07:04
Operator : JU/SJ
Sample : K2063-16DL 10X
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : DB6S4DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM032219MA.M
Quant Title : SVOA CALIBRATION

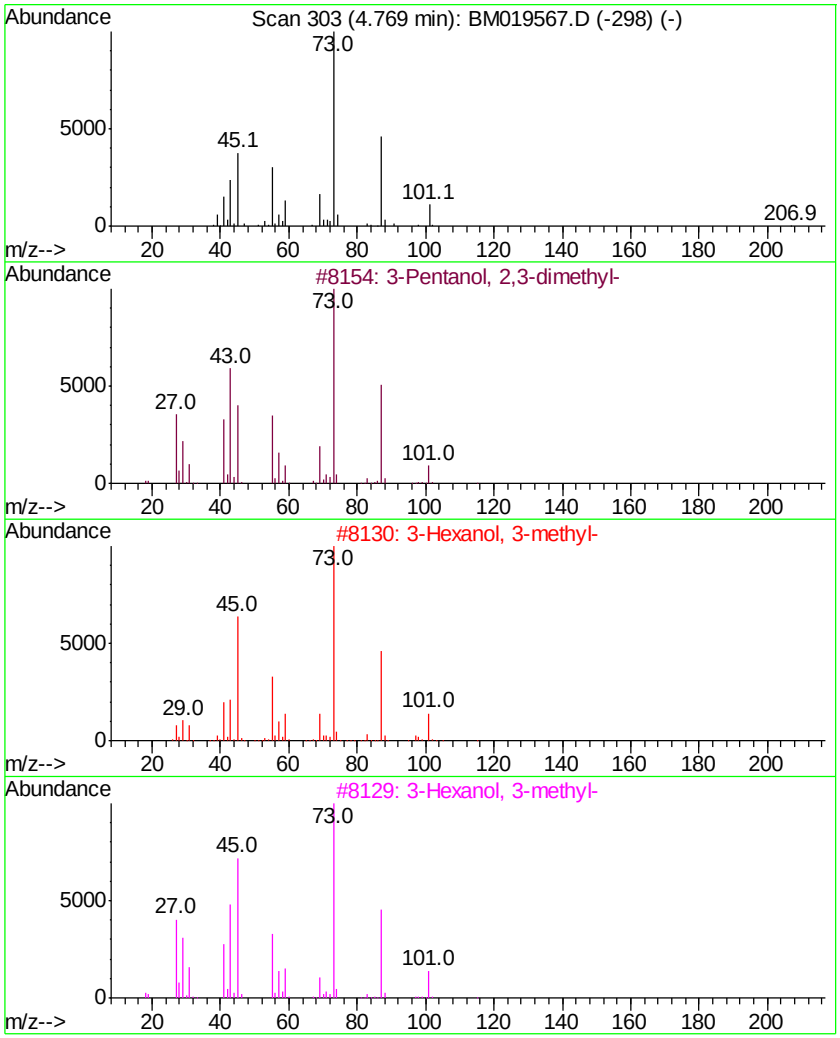
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

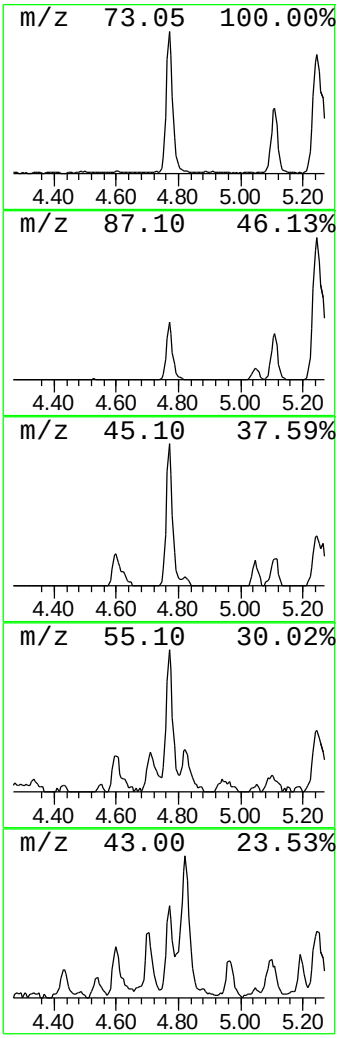
****

Peak Number 6 3-Pentanol, 2,3-dimethyl- Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.77	3.65 ng/ul	194973	1,4-Dichlorobenzene-d4	7.57

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Pentanol, 2,3-dimethyl-	116	C7H16O	000595-41-5	83
2	3-Hexanol, 3-methyl-	116	C7H16O	000597-96-6	74
3	3-Hexanol, 3-methyl-	116	C7H16O	000597-96-6	74
4	3-Hexanol, 3-methyl-	116	C7H16O	000597-96-6	74
5	3-Pentanol, 2,3-dimethyl-	116	C7H16O	000595-41-5	64





Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM032819\
Data File : BM019567.D
Acq On : 29 Mar 2019 07:04
Operator : JU/SJ
Sample : K2063-16DL 10X
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : DB6S4DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM032219MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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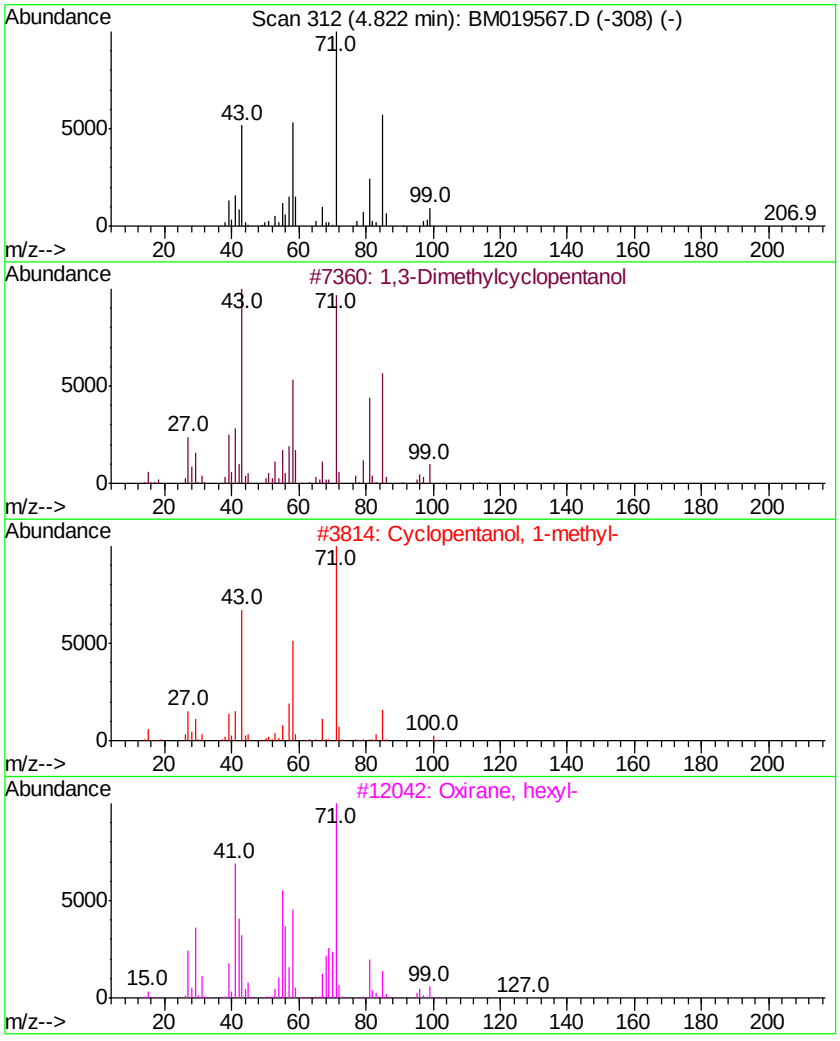
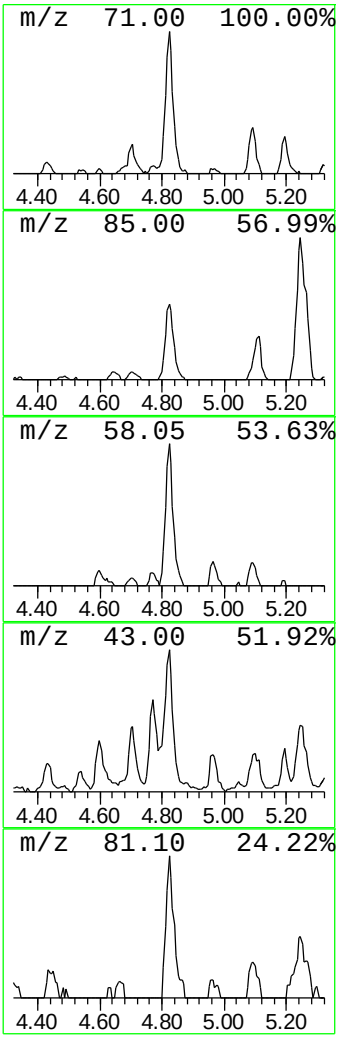
Peak Number 7 1,3-Dimethylcyclopentanol Concentration Rank 28

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.82	3.30 ng/ul	176165	1,4-Dichlorobenzene-d4	7.57

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,3-Dimethylcyclopentanol	114	C7H14O	019550-46-0	64
2	Cyclopentanol, 1-methyl-	100	C6H12O	001462-03-9	42
3	Oxirane, hexyl-	128	C8H16O	002984-50-1	36
4	Thiazole	85	C3H3NS	000288-47-1	35
5	2-Hexanone, 5-methyl-	114	C7H14O	000110-12-3	25





Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM032819\
Data File : BM019567.D
Acq On : 29 Mar 2019 07:04
Operator : JU/SJ
Sample : K2063-16DL 10X
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : DB6S4DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM032219MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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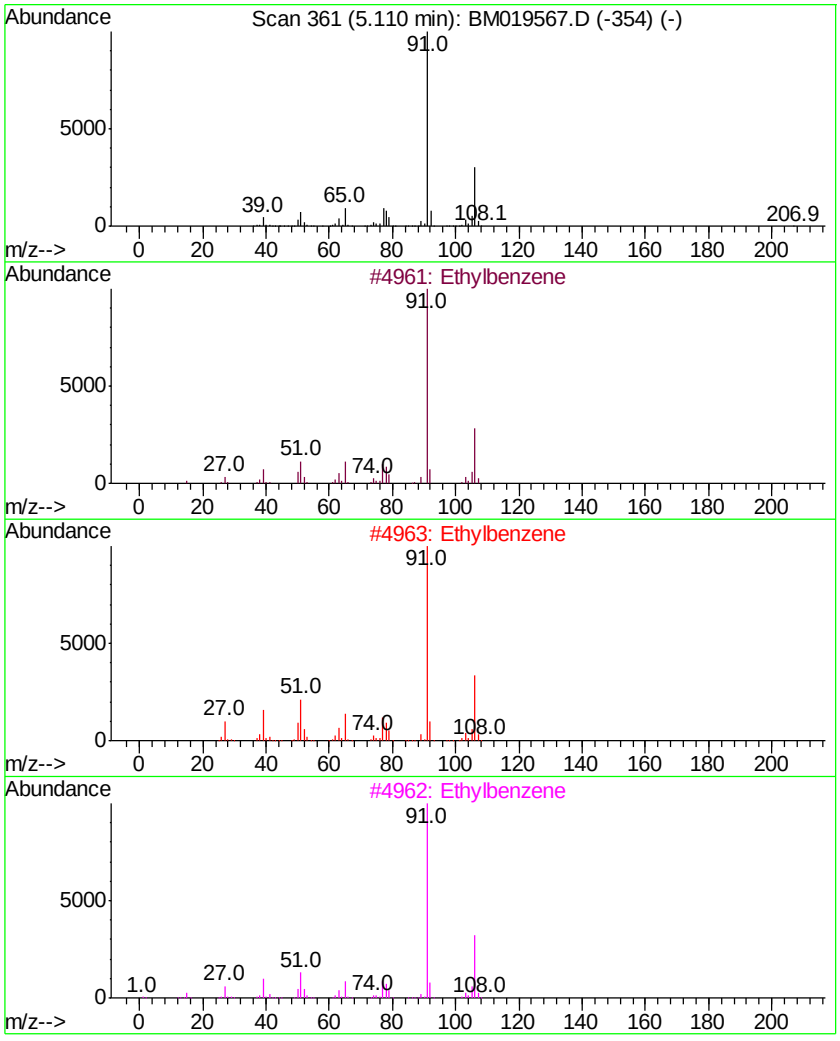
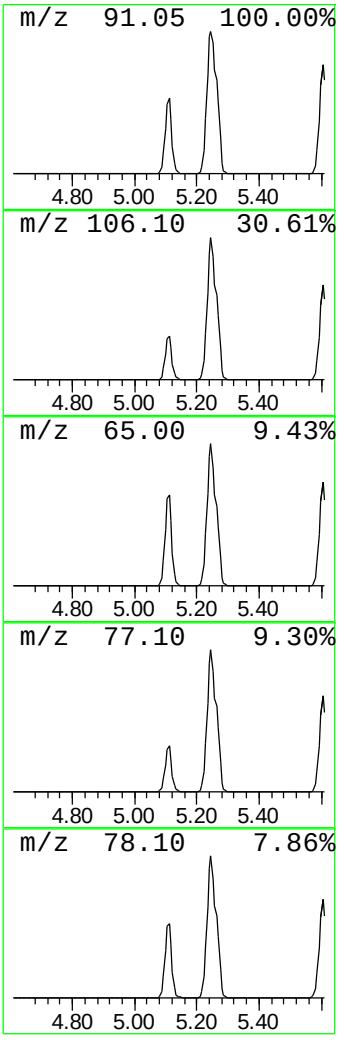
Peak Number 8 1,3-Cyclopentadiene, 5-(1-m... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.11	273.20 ng/ul	14581400	1,4-Dichlorobenzene-d4	7.57

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethylbenzene	106	C8H10	000100-41-4	91
2	Ethylbenzene	106	C8H10	000100-41-4	91
3	Ethylbenzene	106	C8H10	000100-41-4	91
4	Ethylbenzene	106	C8H10	000100-41-4	90
5	1,3-Cyclopentadiene, 5-(1-methyl...	106	C8H10	002175-91-9	80





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Acq On : 29 Mar 2019 07:04
Operator : JU/SJ
Sample : K2063-16DL 10X
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : DB6S4DL

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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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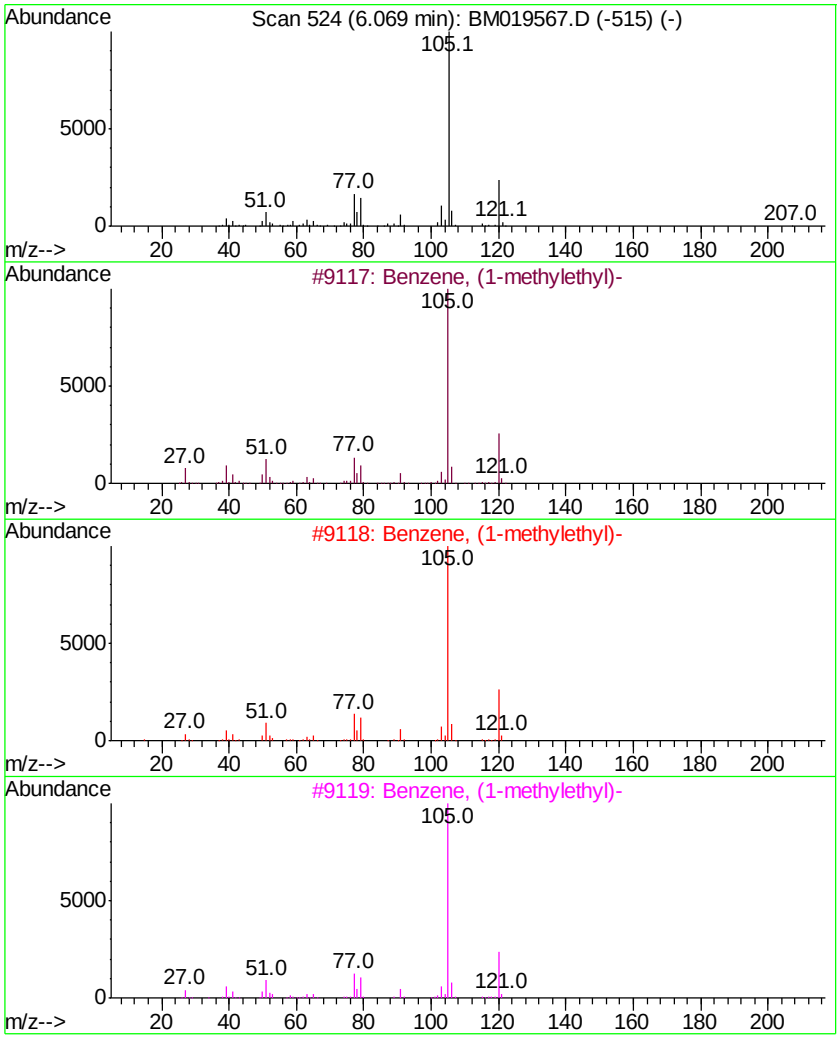
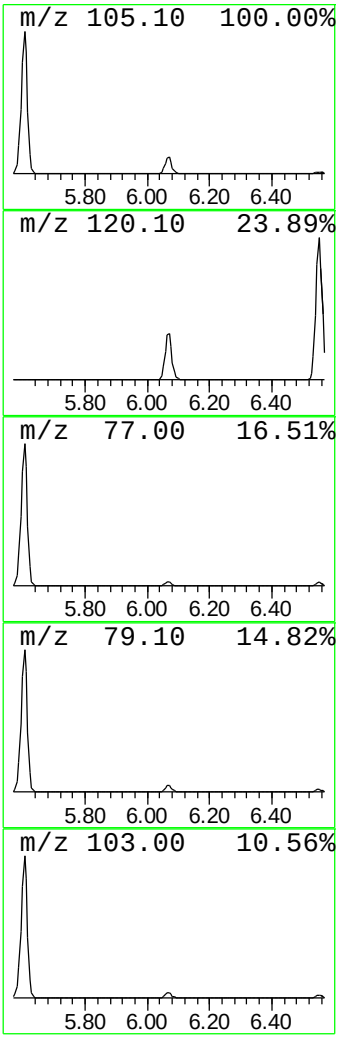
Peak Number 11 Benzene, (1-methylethyl)- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.07	11.48 ng/ul	612490	1,4-Dichlorobenzene-d4	7.57

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, (1-methylethyl)-	120	C9H12	000098-82-8	91
2	Benzene, (1-methylethyl)-	120	C9H12	000098-82-8	91
3	Benzene, (1-methylethyl)-	120	C9H12	000098-82-8	91
4	Benzene, (1-methylethyl)-	120	C9H12	000098-82-8	91
5	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91





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Sample : K2063-16DL 10X
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : DB6S4DL

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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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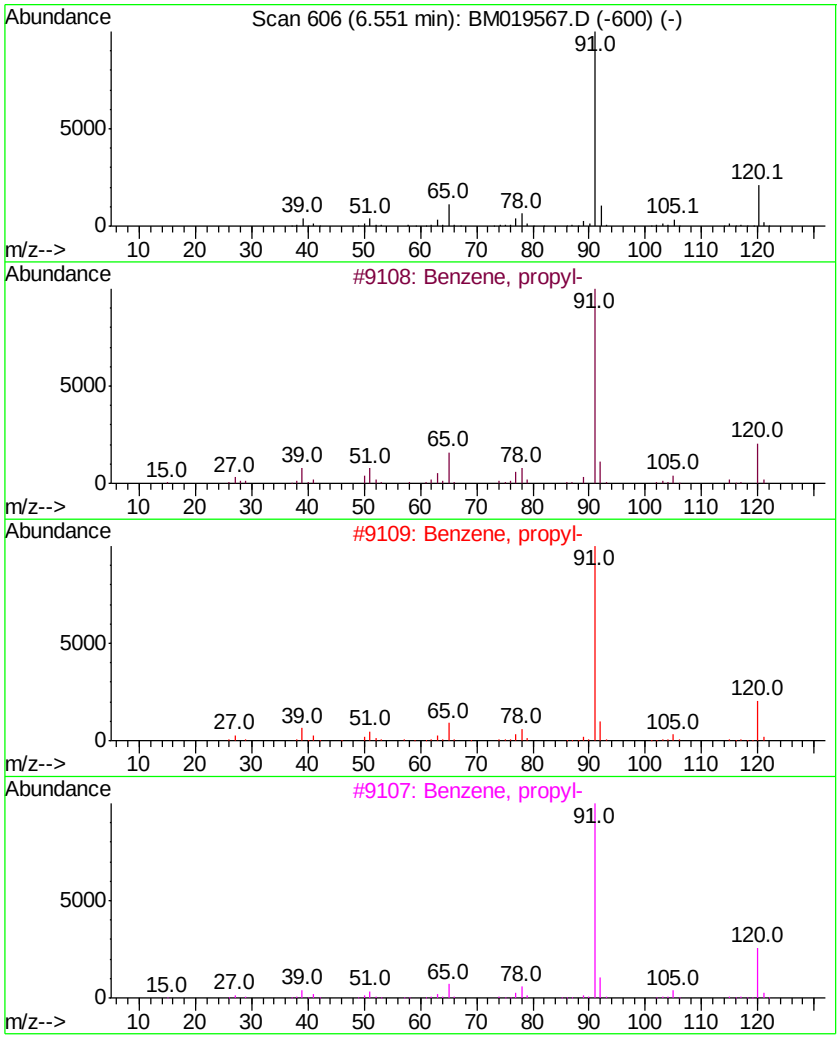
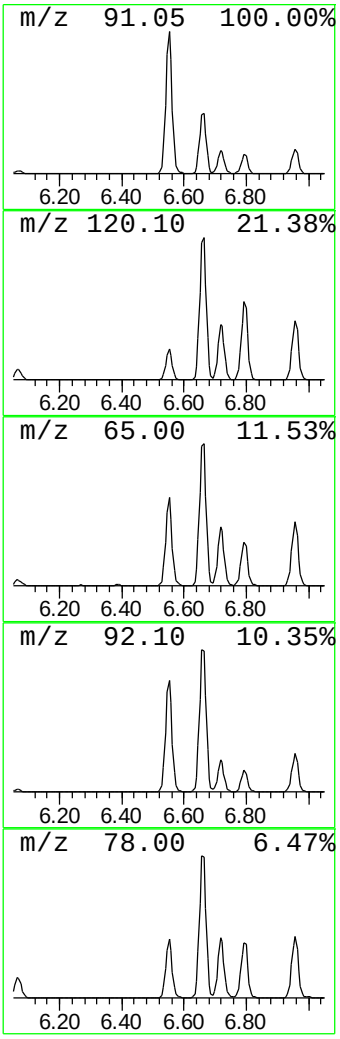
Peak Number 12 N-Benzyl-2-phenethylamine Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.55	26.37 ng/ul	1407400	1,4-Dichlorobenzene-d4	7.57

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, propyl-	120	C9H12	000103-65-1	91
2	Benzene, propyl-	120	C9H12	000103-65-1	90
3	Benzene, propyl-	120	C9H12	000103-65-1	90
4	N-Benzyl-2-phenethylamine	211	C15H17N	003647-71-0	78
5	Phenethylamine, N-benzyl-p-chloro-	245	C15H16ClN	013622-43-0	78





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ClientSampleId : DB6S4DL

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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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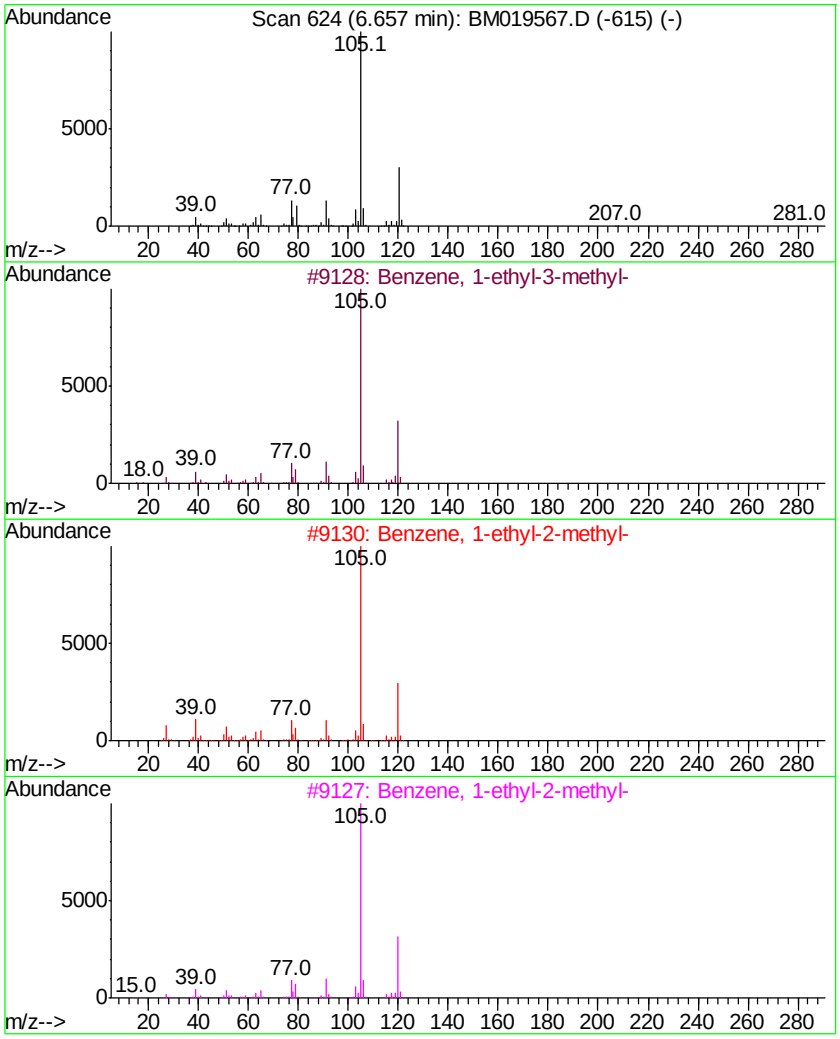
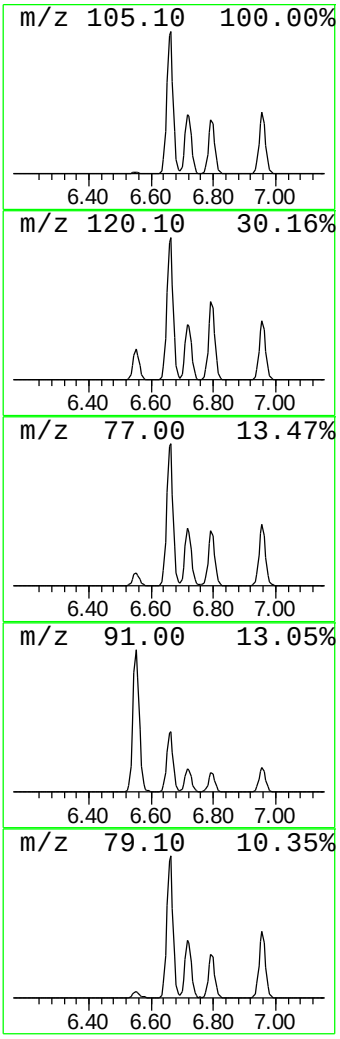
Peak Number 13 Benzene, 1-ethyl-3-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.66	113.33 ng/ul	6048850	1,4-Dichlorobenzene-d4	7.57

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	95
2	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
3	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
4	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	94
5	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91





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ClientSampleId : DB6S4DL

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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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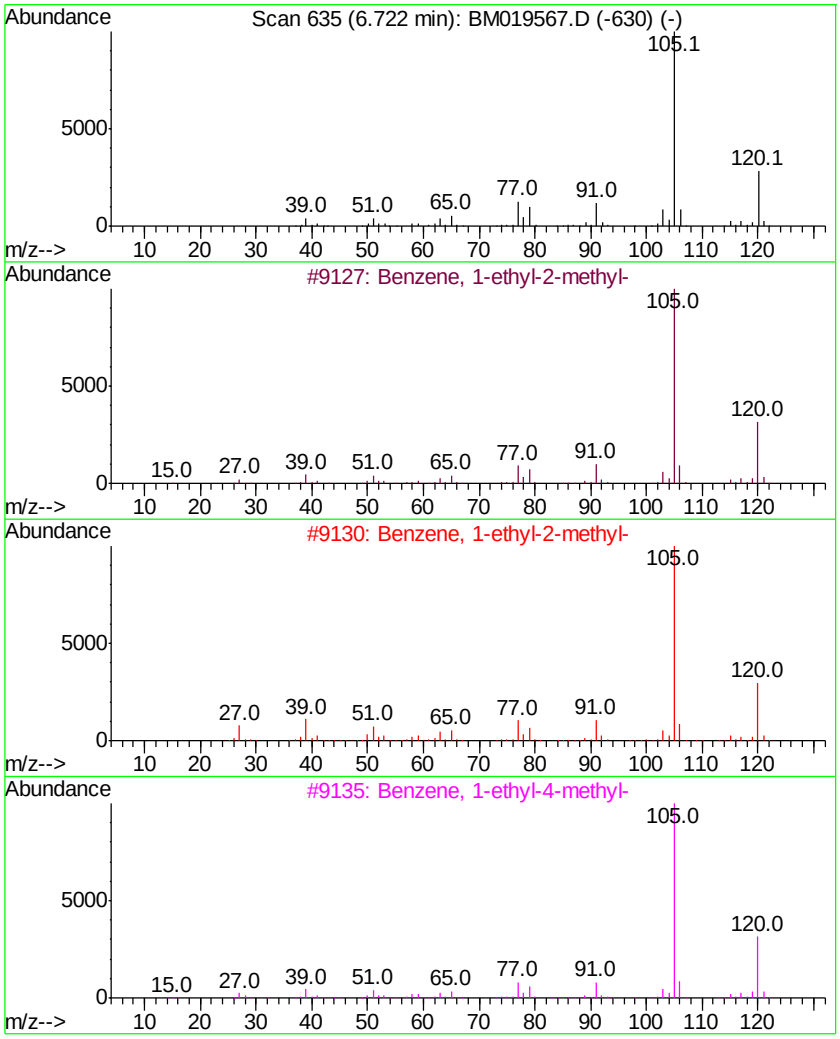
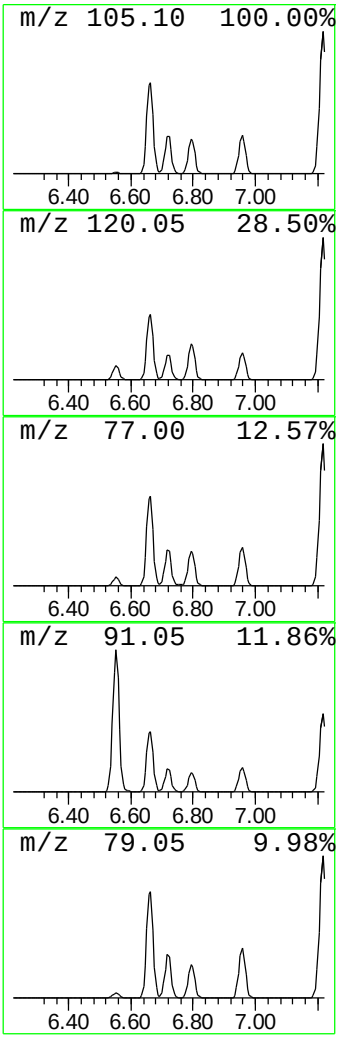
Peak Number 14 Benzene, 1-ethyl-2-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.72	45.54 ng/ul	2430560	1,4-Dichlorobenzene-d4	7.57

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
2	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
3	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
4	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
5	Benzene, 1,3,5-trimethyl-	120	C9H12	000108-67-8	91





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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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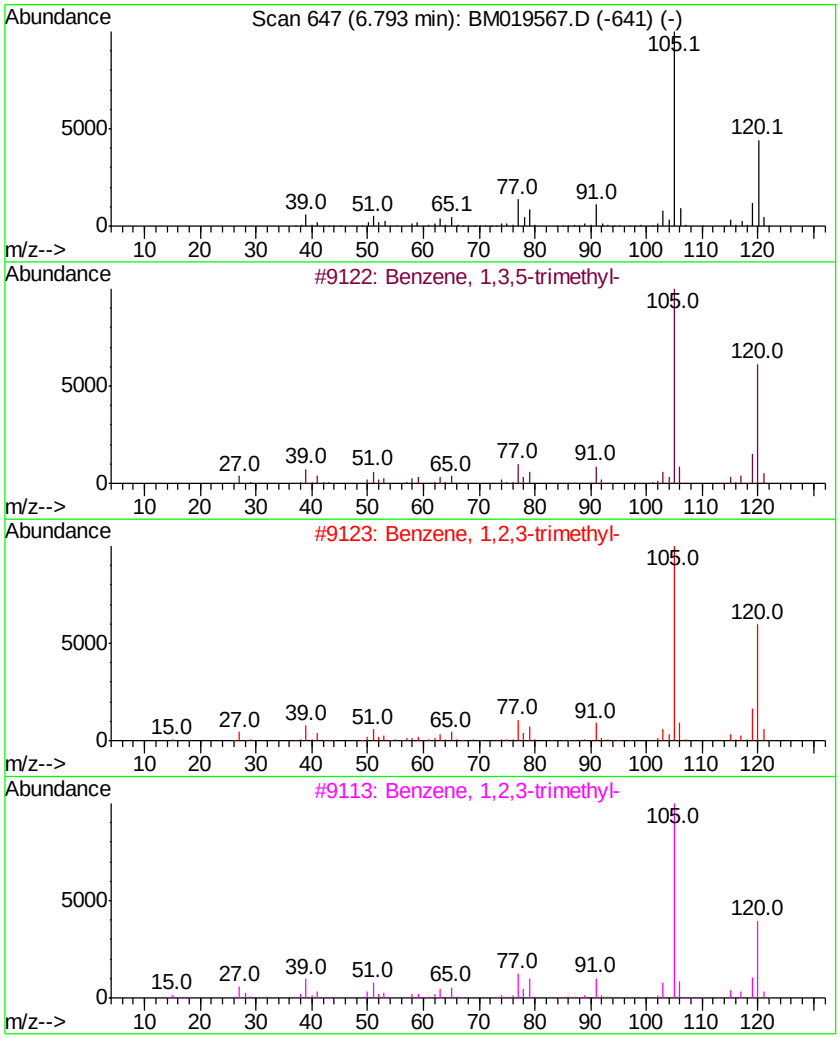
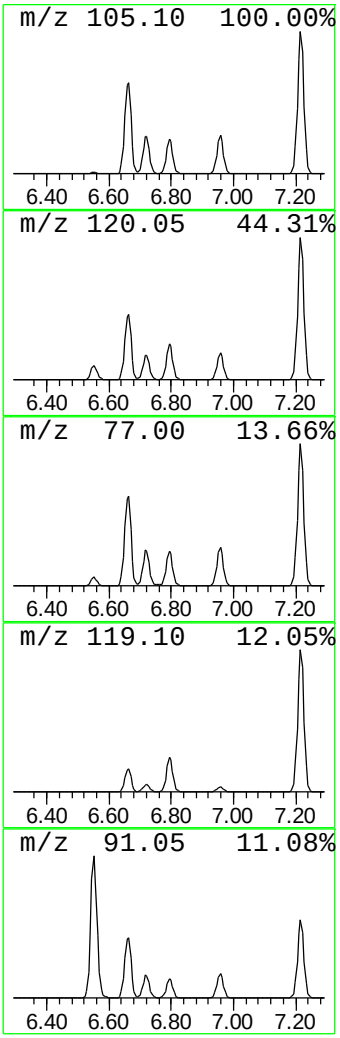
Peak Number 15 Benzene, 1,3,5-trimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.79	46.97 ng/ul	2506950	1,4-Dichlorobenzene-d4	7.57

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,3,5-trimethyl-	120	C9H12	000108-67-8	97
2	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	97
3	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
4	Benzene, 1,3,5-trimethyl-	120	C9H12	000108-67-8	94
5	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94





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Sample : K2063-16DL 10X
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ALS Vial : 29 Sample Multiplier: 1

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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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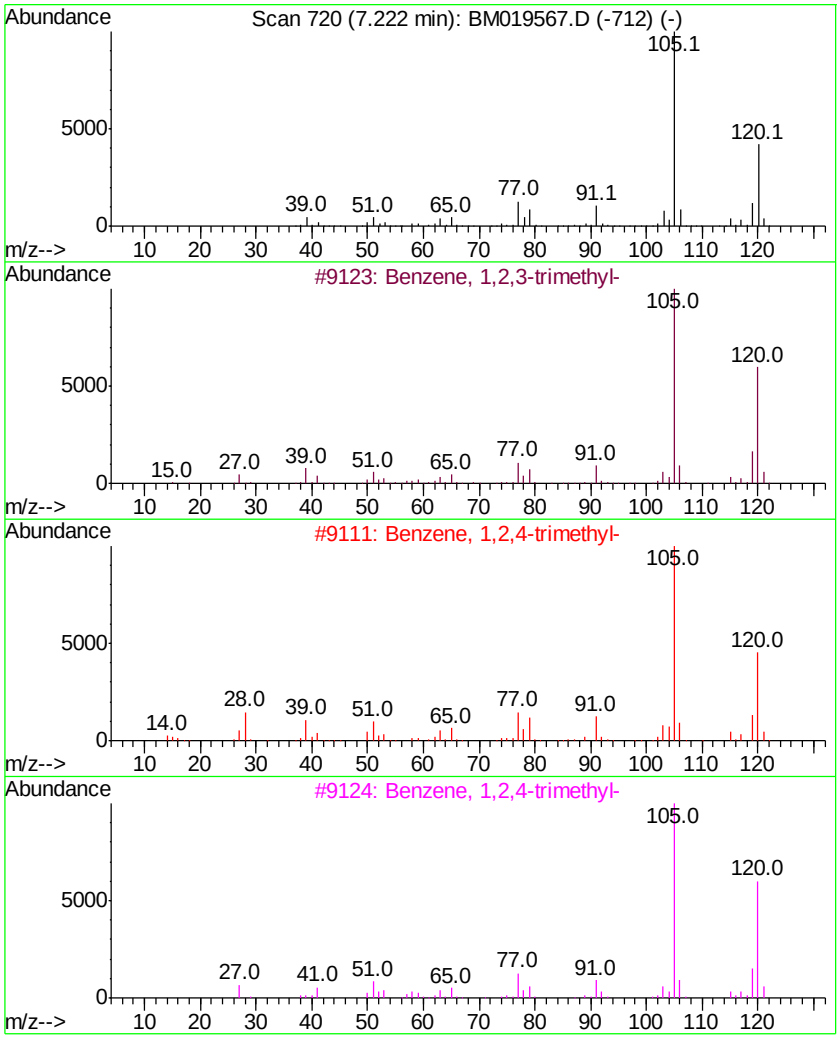
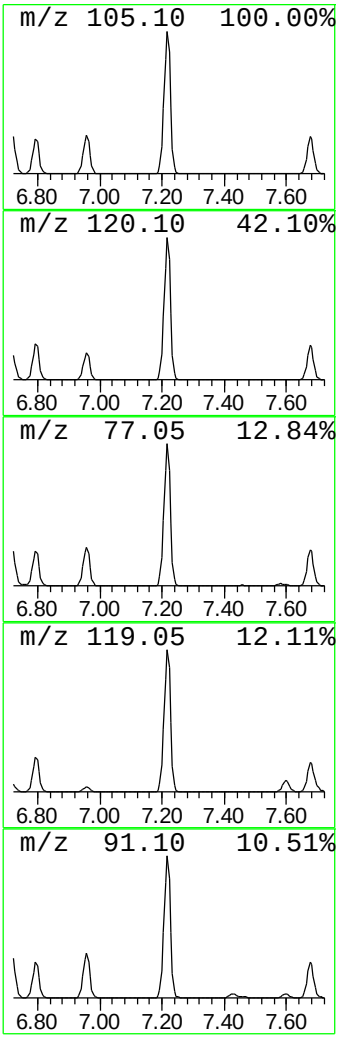
Peak Number 16 Benzene, 1,2,3-trimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.22	189.85 ng/ul	10133100	1,4-Dichlorobenzene-d4	7.57

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	97
2	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94
3	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94
4	Benzene, 1,3,5-trimethyl-	120	C9H12	000108-67-8	91
5	Benzene, 1,3,5-trimethyl-	120	C9H12	000108-67-8	91





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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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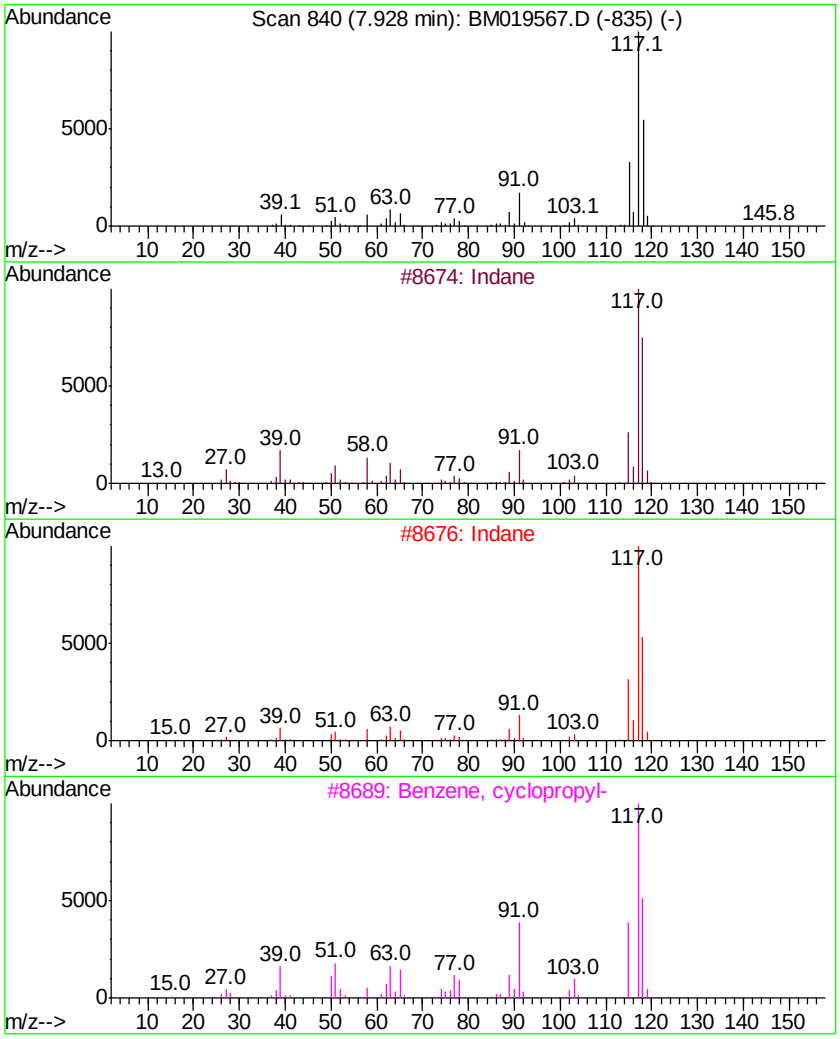
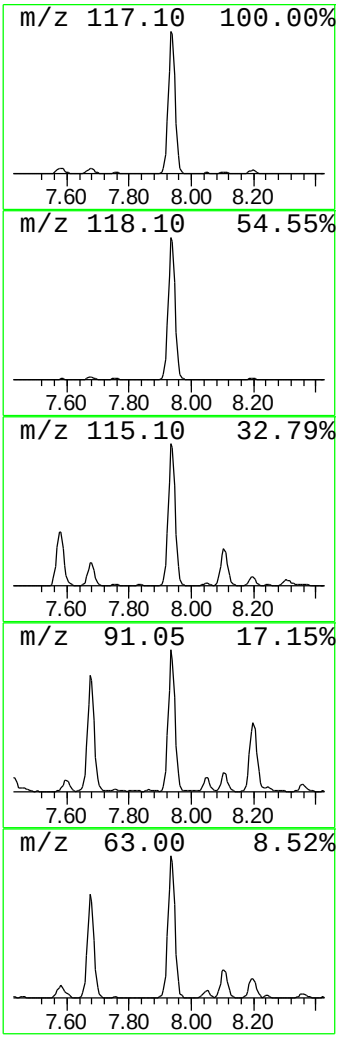
Peak Number 18 Indane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.93	42.91 ng/ul	2290430	1,4-Dichlorobenzene-d4	7.57

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane	118	C9H10	000496-11-7	91
2	Indane	118	C9H10	000496-11-7	91
3	Benzene, cyclopropyl-	118	C9H10	000873-49-4	72
4	Deltacyclene	118	C9H10	007785-10-6	72
5	Benzene, 1,1'-(1,5-hexadiene-1,6...	234	C18H18	004439-45-6	59





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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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Peak Number 19 Benzene, 1,3-diethyl- Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.05	3.86 ng/ul	205762	1,4-Dichlorobenzene-d4	7.57

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	94
2	Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	94
3	Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	93
4	Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	93
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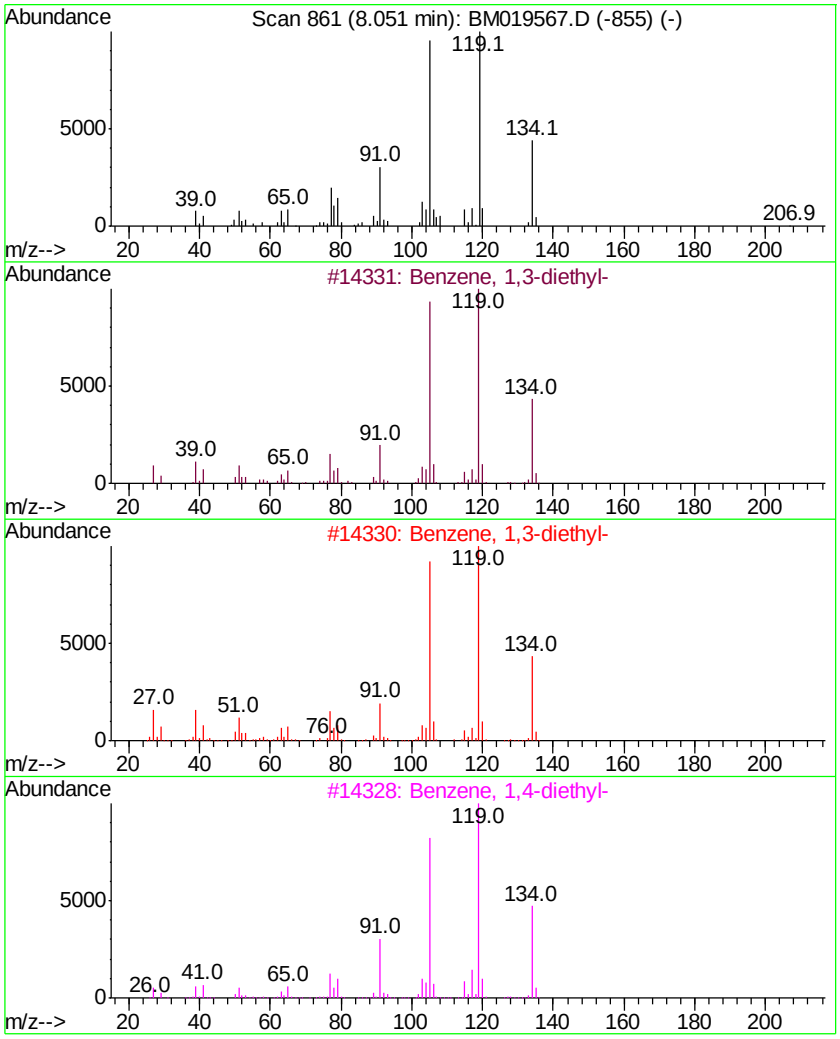

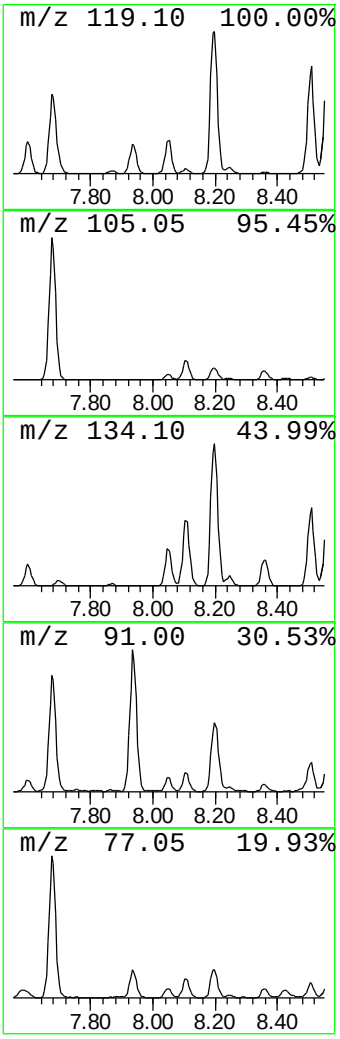

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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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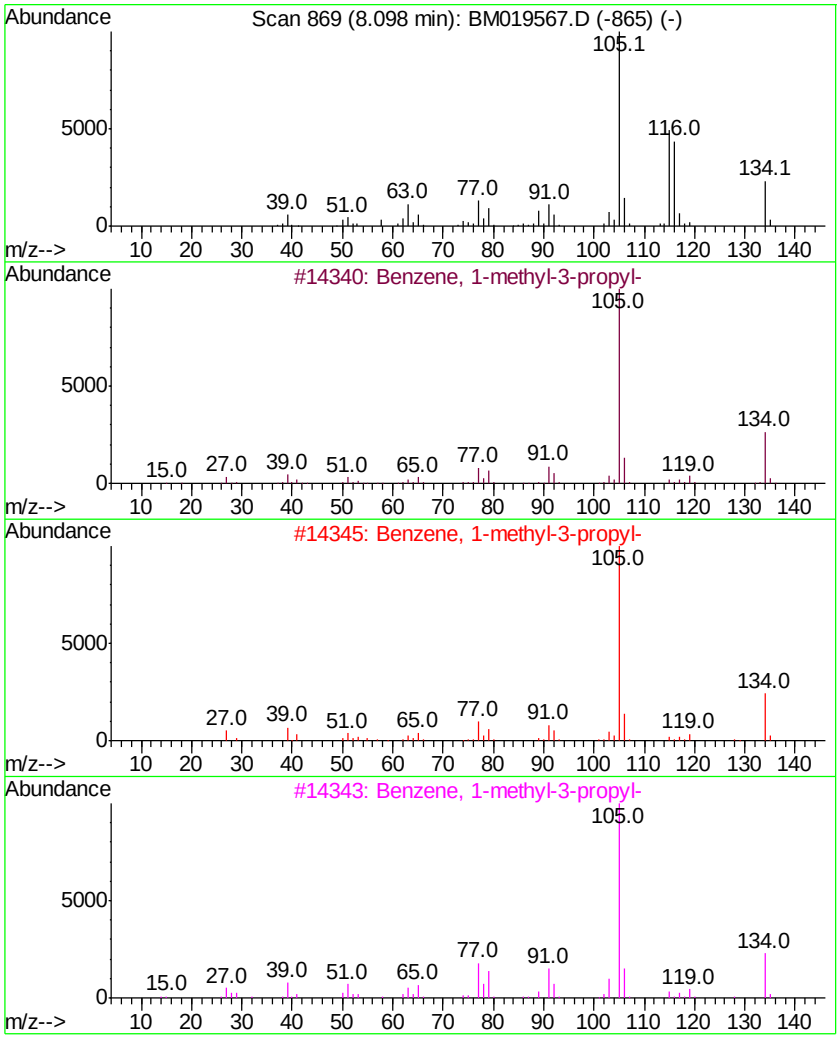
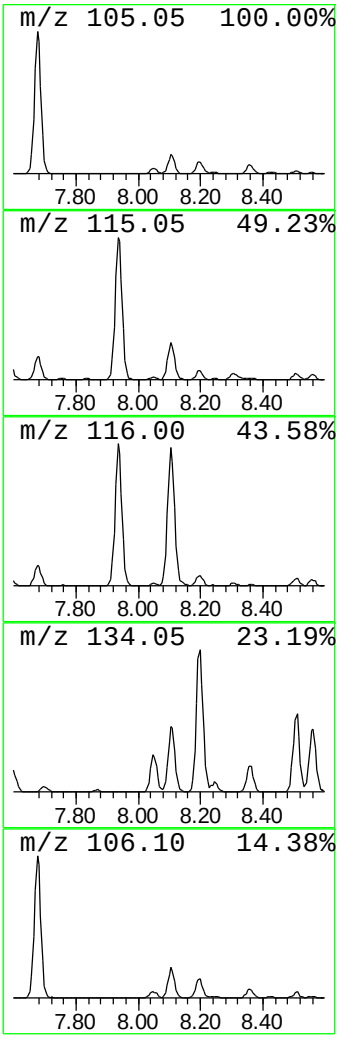
Peak Number 20 unknown-01 Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.10	10.07 ng/ul	537667	1,4-Dichlorobenzene-d4	7.57

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-methyl-3-propyl-	134	C10H14	001074-43-7	46
2	Benzene, 1-methyl-3-propyl-	134	C10H14	001074-43-7	46
3	Benzene, 1-methyl-3-propyl-	134	C10H14	001074-43-7	42
4	Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	42
5	Benzene, 1-methyl-2-propyl-	134	C10H14	001074-17-5	42





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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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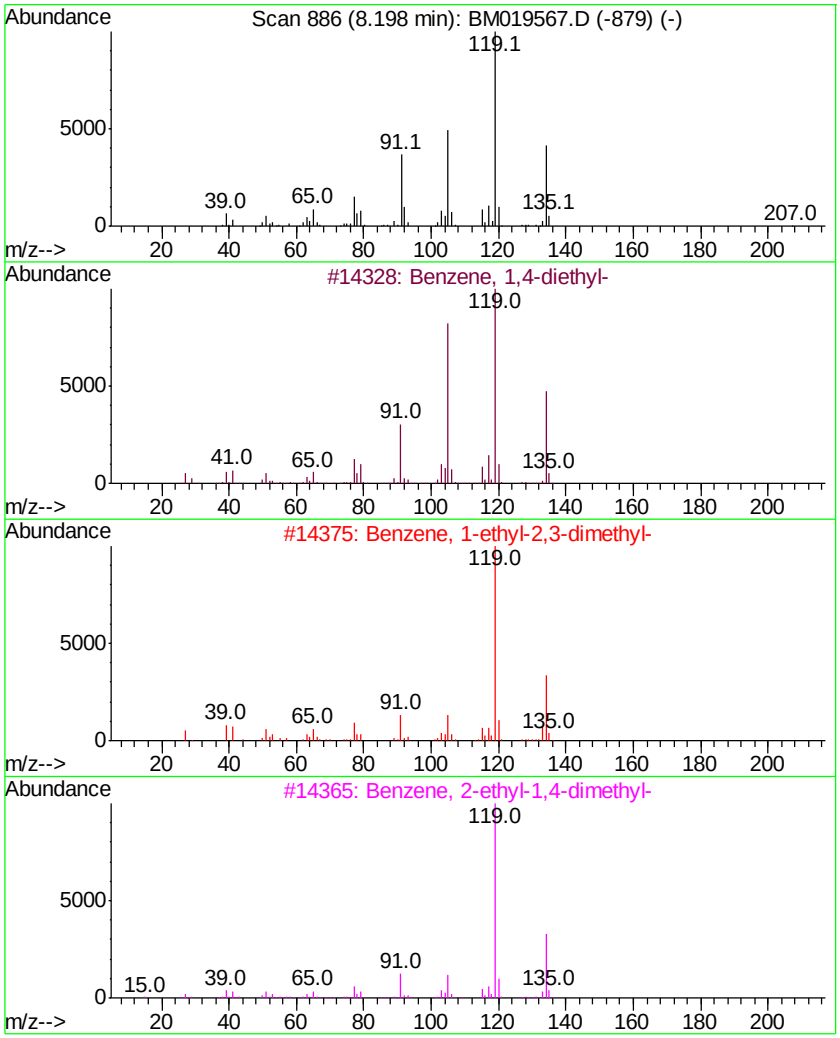
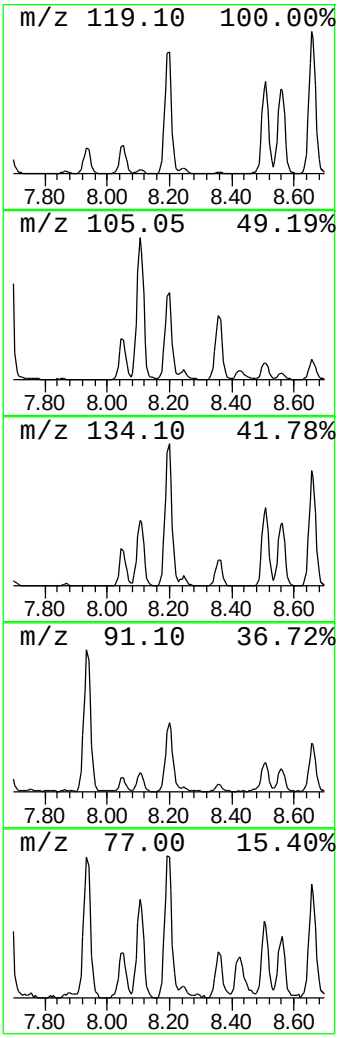
Peak Number 21 Benzene, 1,4-diethyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.20	13.20 ng/ul	704470	1,4-Dichlorobenzene-d4	7.57

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	92
2	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	91
3	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	91
4	Benzene, 1-methyl-2-(1-methyleth...	134	C10H14	000527-84-4	91
5	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	91





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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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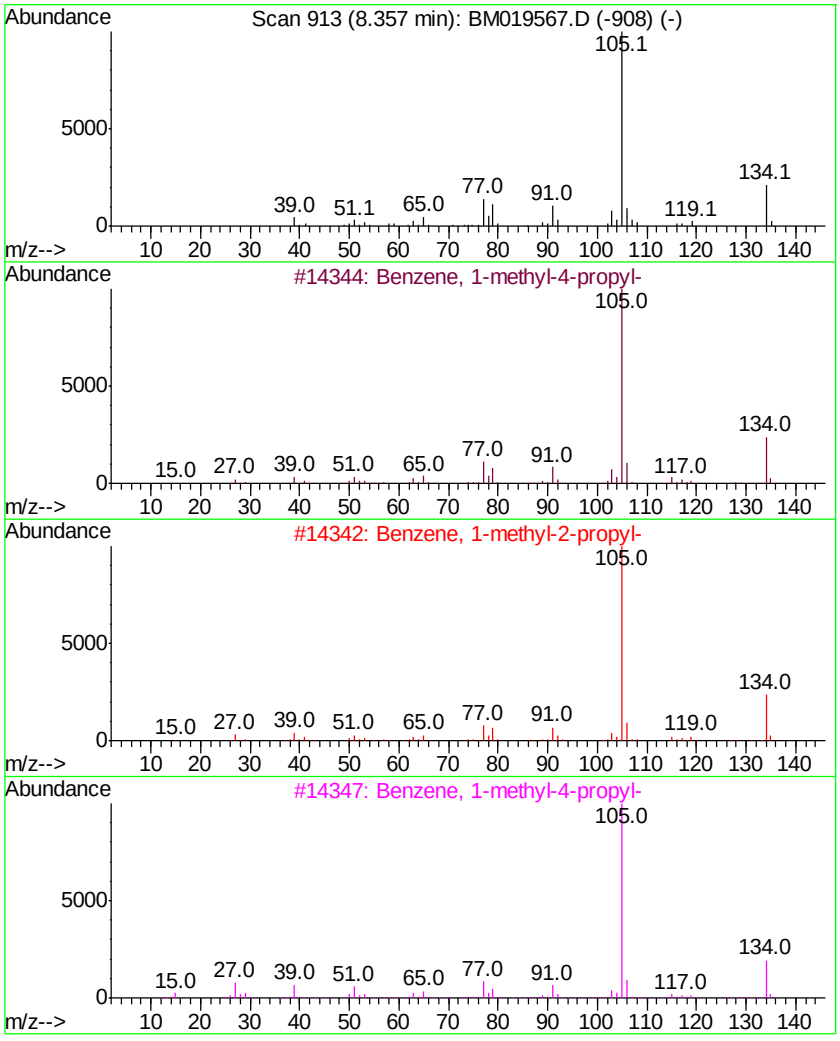
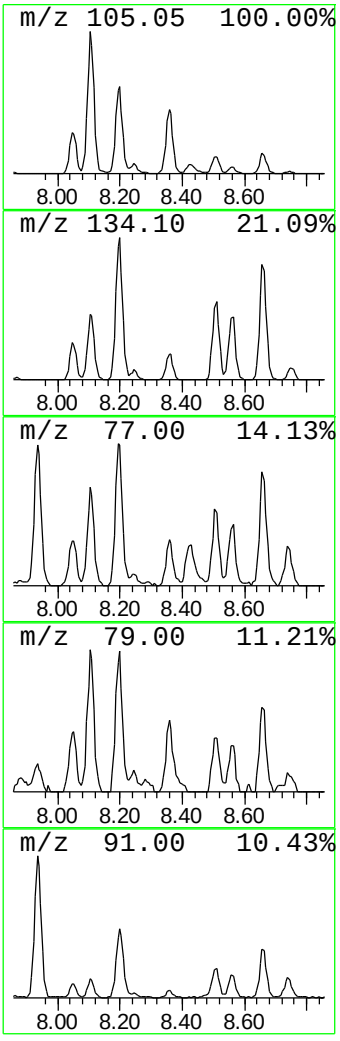
Peak Number 22 Benzene, 1-methyl-4-propyl- Concentration Rank 30

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.36	2.80 ng/ul	149368	1,4-Dichlorobenzene-d4	7.57

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	94
2	Benzene, 1-methyl-2-propyl-	134	C10H14	001074-17-5	91
3	Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	91
4	Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	91
5	Benzene, (1-methylpropyl)-	134	C10H14	000135-98-8	90





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Data File : BM019567.D
Acq On : 29 Mar 2019 07:04
Operator : JU/SJ
Sample : K2063-16DL 10X
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : DB6S4DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM032219MA.M
Quant Title : SVOA CALIBRATION

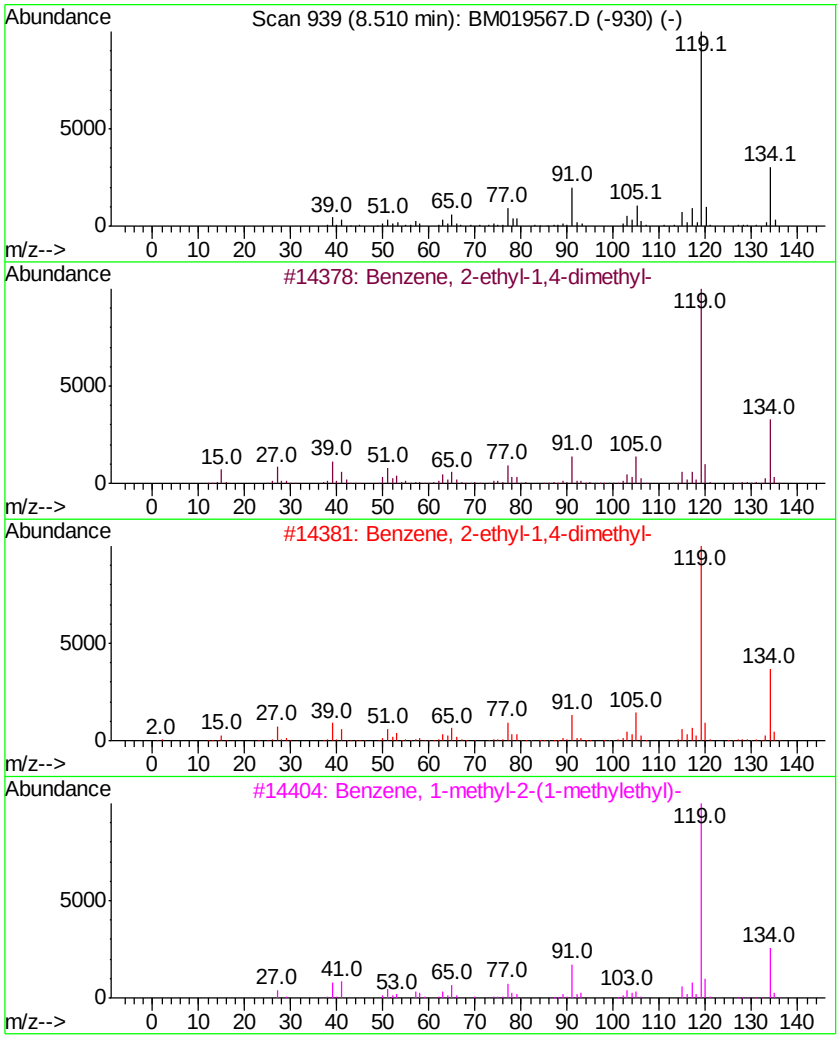
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

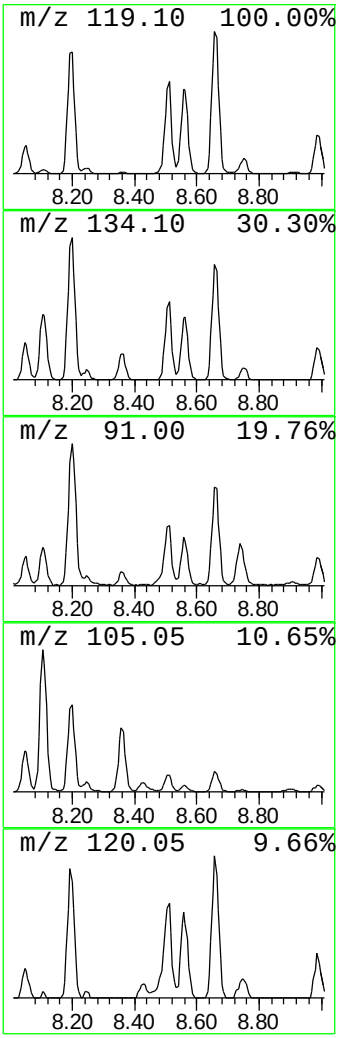
****

Peak Number 23 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.51	7.68 ng/ul	409808	1,4-Dichlorobenzene-d4	7.57

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	95
2	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	95
3	Benzene, 1-methyl-2-(1-methyleth...	134	C10H14	000527-84-4	95
4	Benzene, 1-methyl-2-(1-methyleth...	134	C10H14	000527-84-4	95
5	Benzene, 1-methyl-4-(1-methyleth...	134	C10H14	000099-87-6	95





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Data File : BM019567.D
Acq On : 29 Mar 2019 07:04
Operator : JU/SJ
Sample : K2063-16DL 10X
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : DB6S4DL

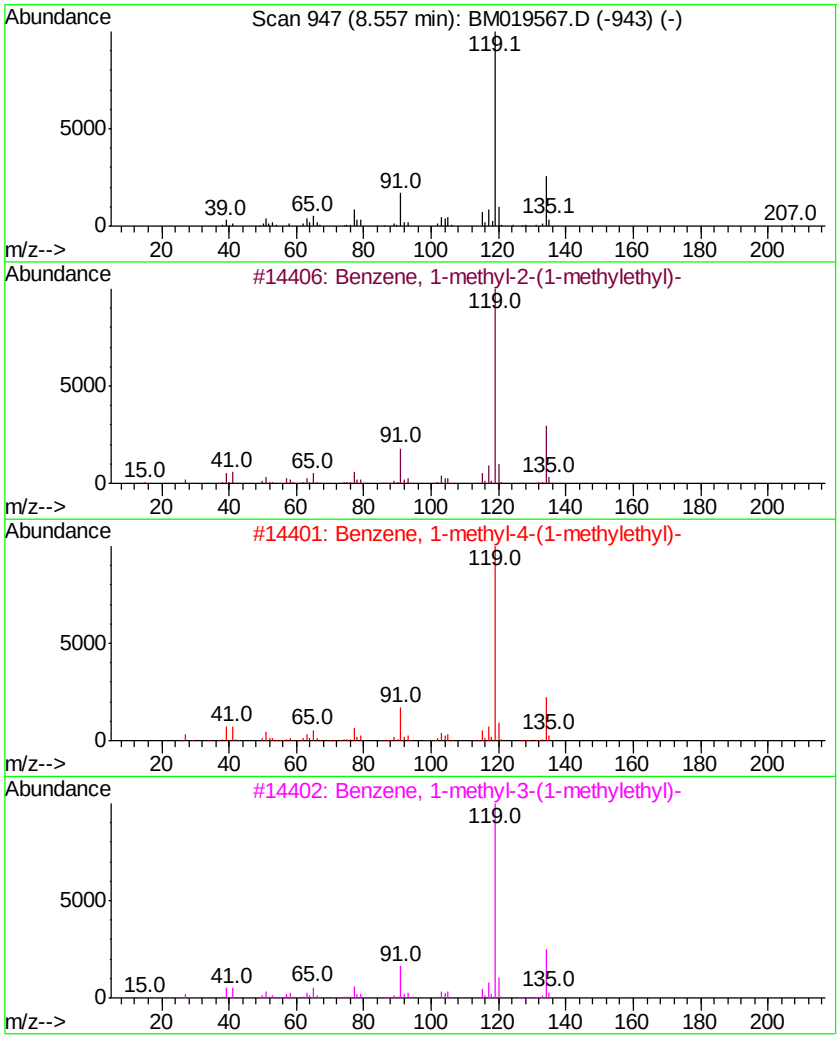
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM032219MA.M
Quant Title : SVOA CALIBRATION

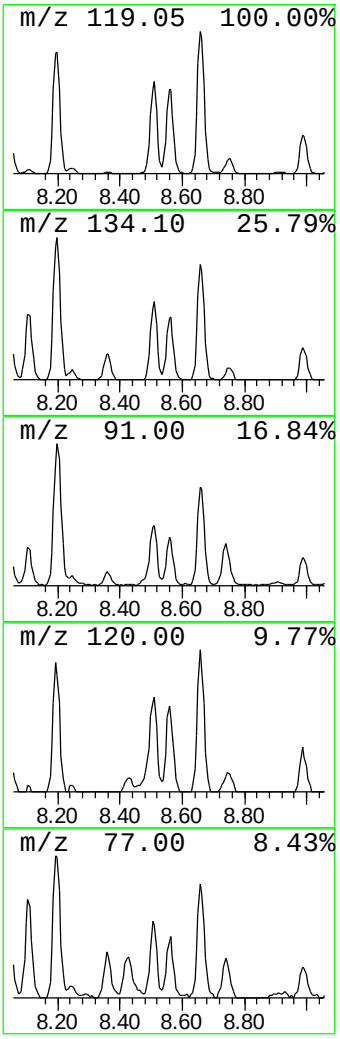
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 24 Benzene, 1-methyl-2-(1-meth... Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.56	6.69 ng/ul	357047	1,4-Dichlorobenzene-d4	7.57

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-methyl-2-(1-methyleth...	134	C10H14	000527-84-4	97
2	Benzene, 1-methyl-4-(1-methyleth...	134	C10H14	000099-87-6	95
3	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	95
4	Benzene, 1-methyl-2-(1-methyleth...	134	C10H14	000527-84-4	95
5	Benzene, 1-methyl-2-(1-methyleth...	134	C10H14	000527-84-4	95





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Data File : BM019567.D
Acq On : 29 Mar 2019 07:04
Operator : JU/SJ
Sample : K2063-16DL 10X
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : DB6S4DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM032219MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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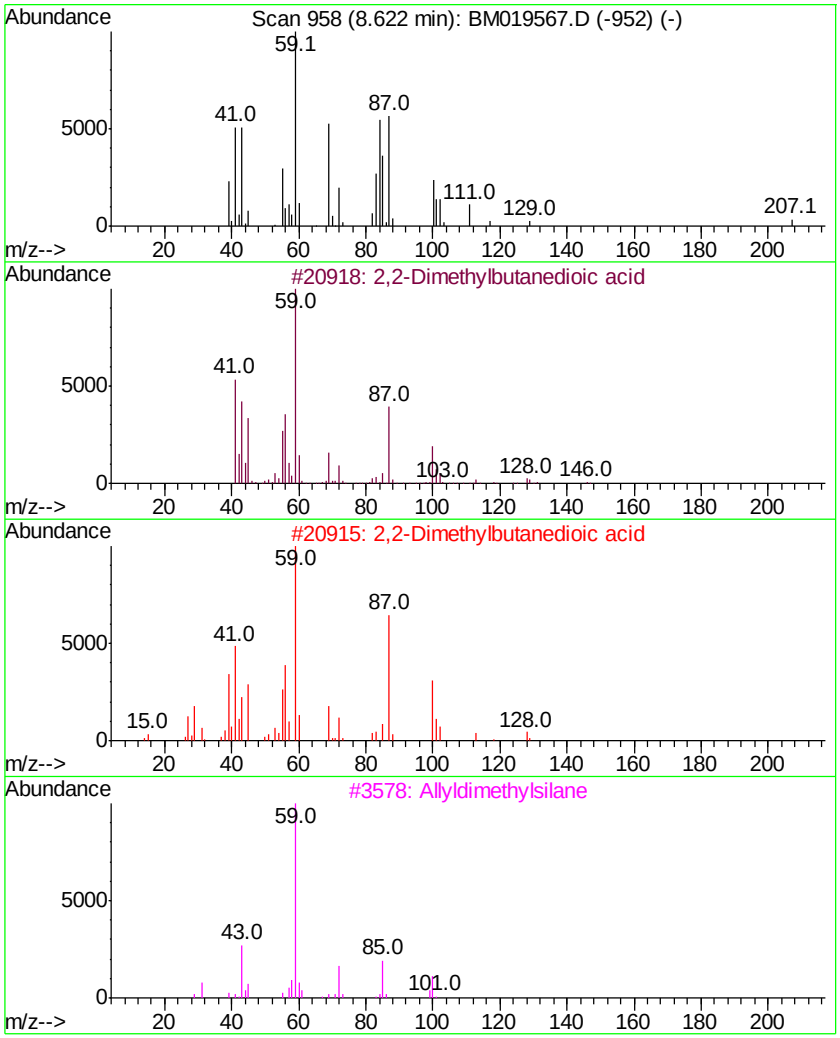
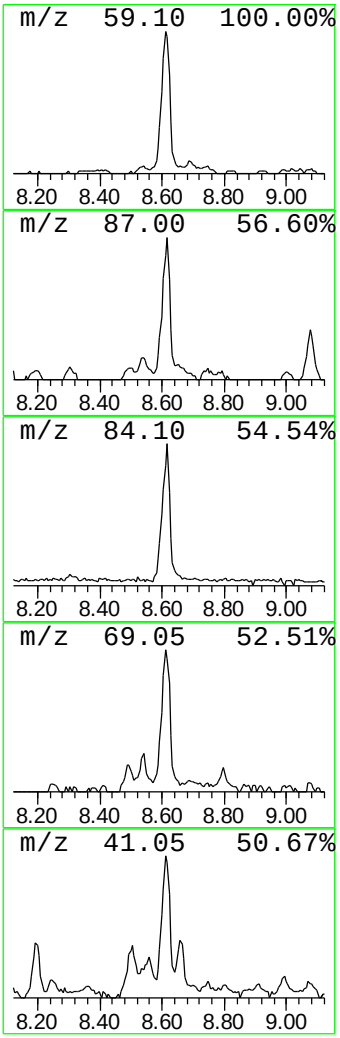
Peak Number 25 unknown-02 Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.62	4.55 ng/ul	242908	1,4-Dichlorobenzene-d4	7.57

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,2-Dimethylbutanedioic acid	146	C6H10O4	000597-43-3	40
2	2,2-Dimethylbutanedioic acid	146	C6H10O4	000597-43-3	38
3	Allyldimethylsilane	100	C5H12Si	003937-30-2	37
4	1,3-Dioxolane-2-propanol, 2-methyl-	146	C7H14O3	029021-98-5	32
5	Butane, 2-ethoxy-2-methyl-	116	C7H16O	000919-94-8	32





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Data File : BM019567.D
Acq On : 29 Mar 2019 07:04
Operator : JU/SJ
Sample : K2063-16DL 10X
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : DB6S4DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM032219MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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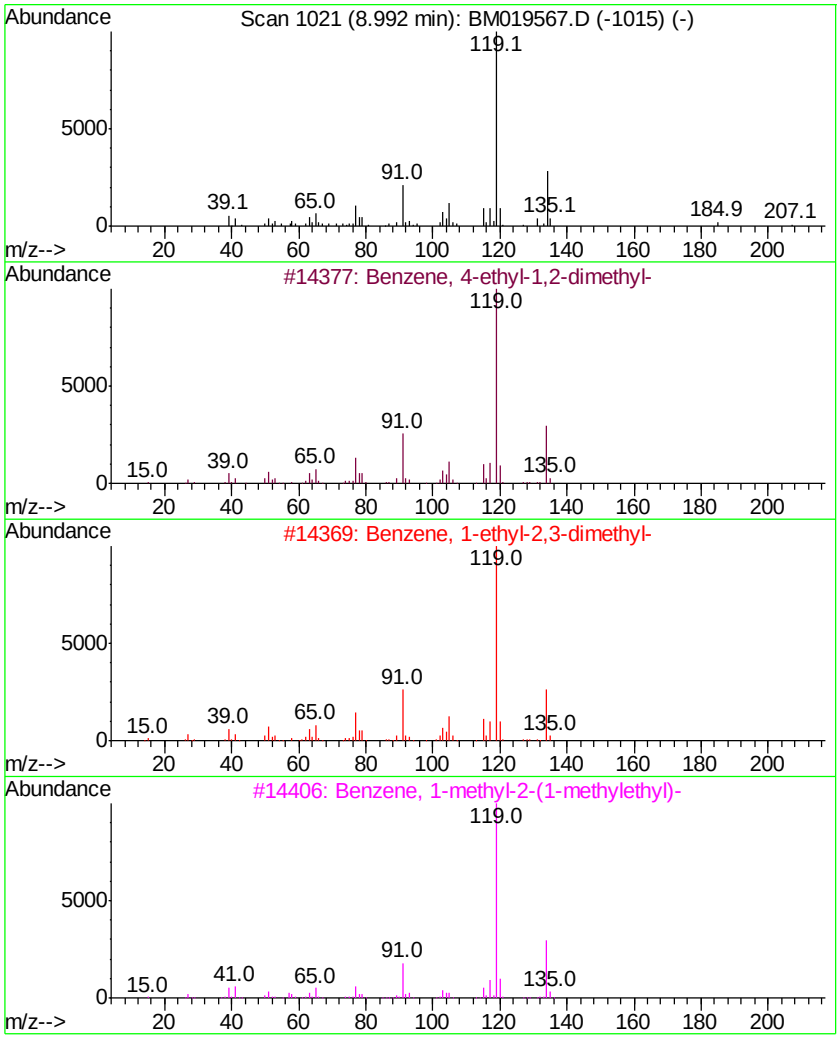
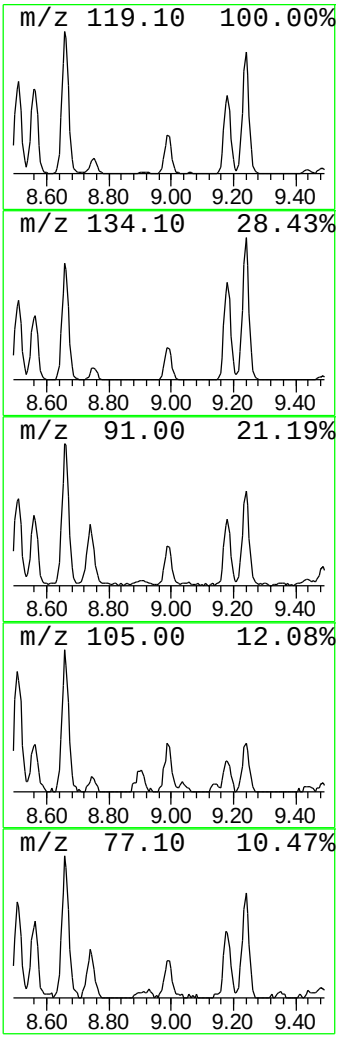
Peak Number 27 Benzene, 4-ethyl-1,2-dimethyl- Concentration Rank 32

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.99	2.47 ng/ul	183276	Naphthalene-d8	10.36

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	96
2	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	96
3	Benzene, 1-methyl-2-(1-methyleth...	134	C10H14	000527-84-4	95
4	Benzene, 1-methyl-2-(1-methyleth...	134	C10H14	000527-84-4	95
5	Benzene, 1-methyl-4-(1-methyleth...	134	C10H14	000099-87-6	95





Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM032819\
Data File : BM019567.D
Acq On : 29 Mar 2019 07:04
Operator : JU/SJ
Sample : K2063-16DL 10X
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : DB6S4DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM032219MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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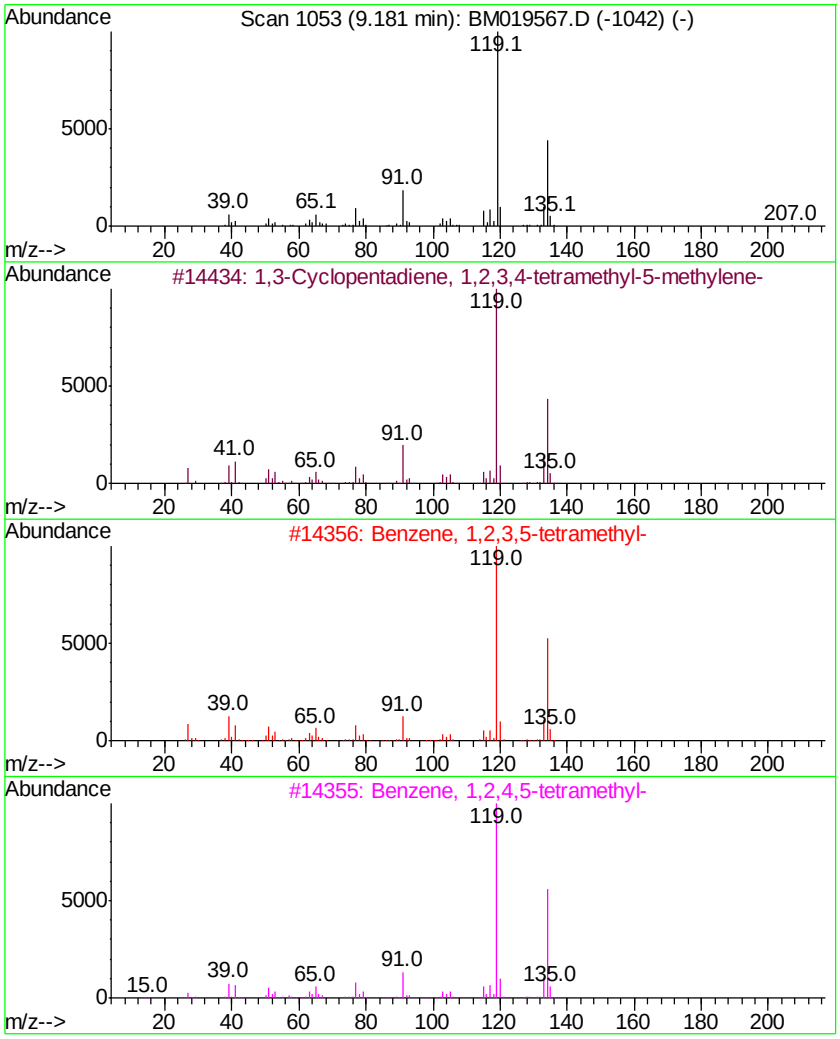
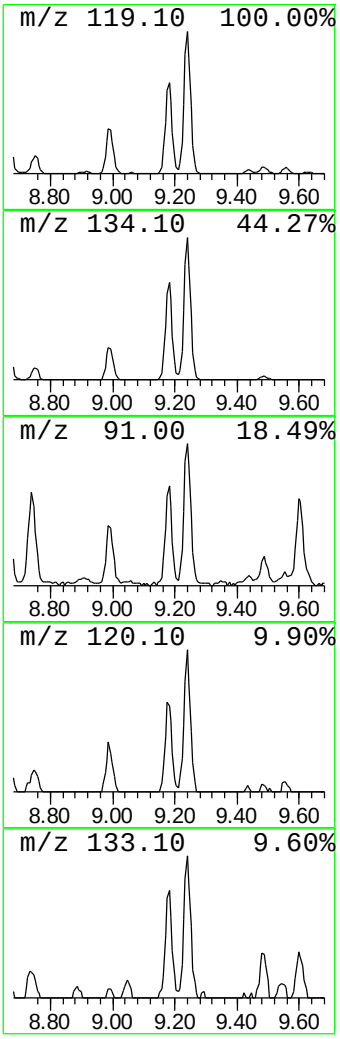
Peak Number 28 1,3-Cyclopentadiene, 1,2,3,... Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.18	4.48 ng/ul	332086	Naphthalene-d8	10.36

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,3-Cyclopentadiene, 1,2,3,4-tet...	134	C10H14	076089-59-3	97
2	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	96
3	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	95
4	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	95
5	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	95





Library Search Compound Report

Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM032819\
Data File : BM019567.D
Acq On : 29 Mar 2019 07:04
Operator : JU/SJ
Sample : K2063-16DL 10X
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : DB6S4DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM032219MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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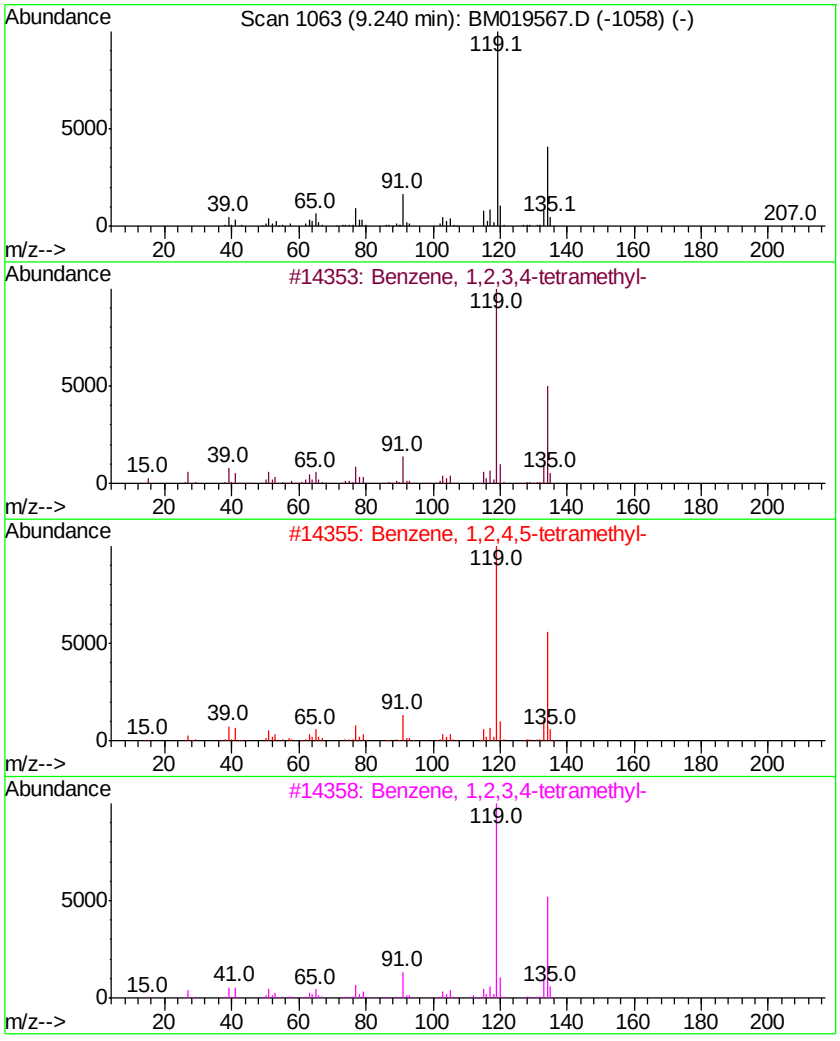
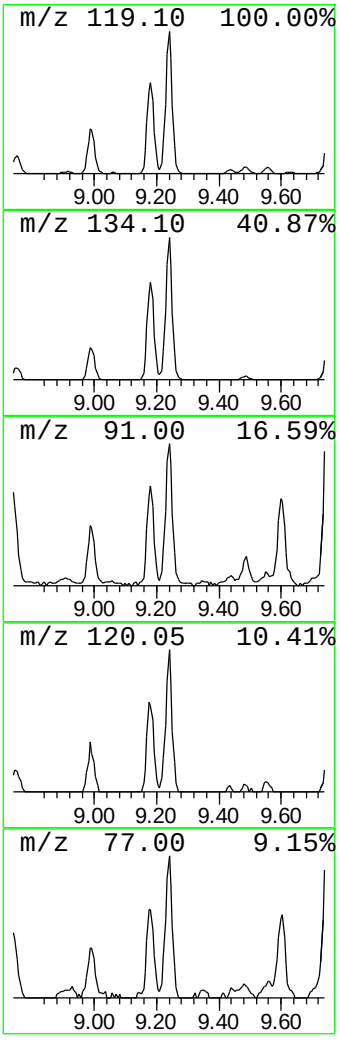
Peak Number 29 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.24	6.26 ng/ul	464460	Naphthalene-d8	10.36

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
2	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	95
3	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	95
4	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	95
5	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	95





Library Search Compound Report

Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM032819\
Data File : BM019567.D
Acq On : 29 Mar 2019 07:04
Operator : JU/SJ
Sample : K2063-16DL 10X
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : DB6S4DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM032219MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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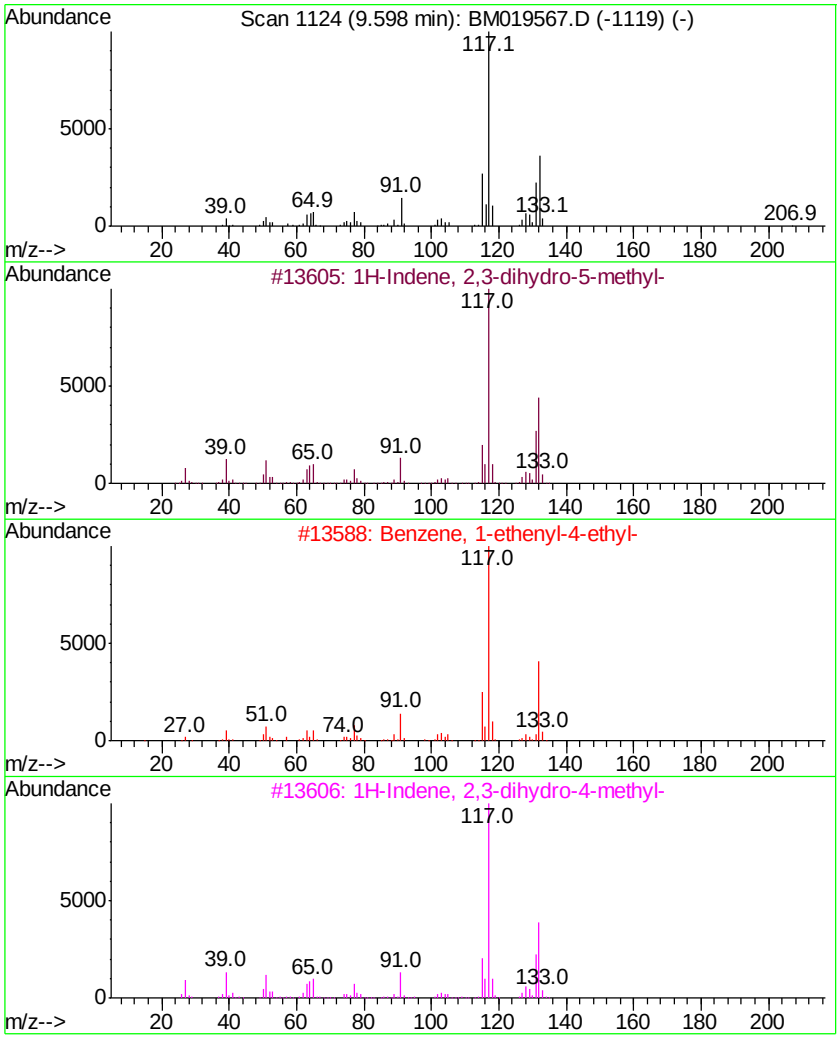
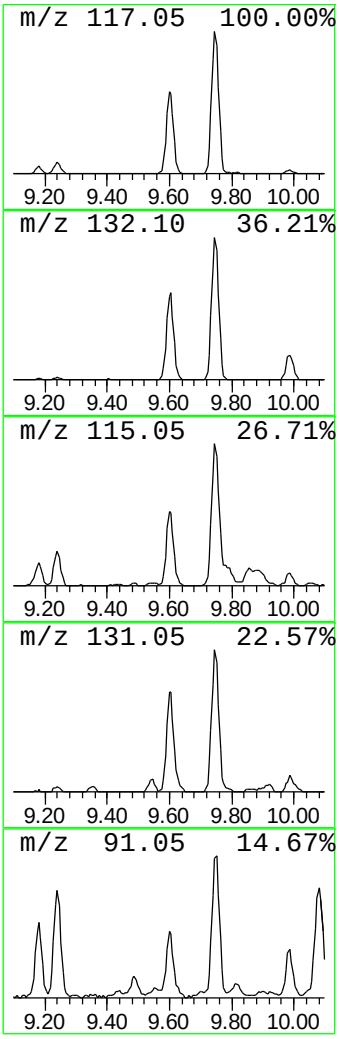
Peak Number 30 1H-Indene, 2,3-dihydro-5-me... Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.60	5.42 ng/ul	402426	Naphthalene-d8	10.36

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	95
2	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	94
3	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	94
4	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	93
5	3-Phenylbut-1-ene	132	C10H12	000934-10-1	91





Library Search Compound Report

Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM032819\
Data File : BM019567.D
Acq On : 29 Mar 2019 07:04
Operator : JU/SJ
Sample : K2063-16DL 10X
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : DB6S4DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM032219MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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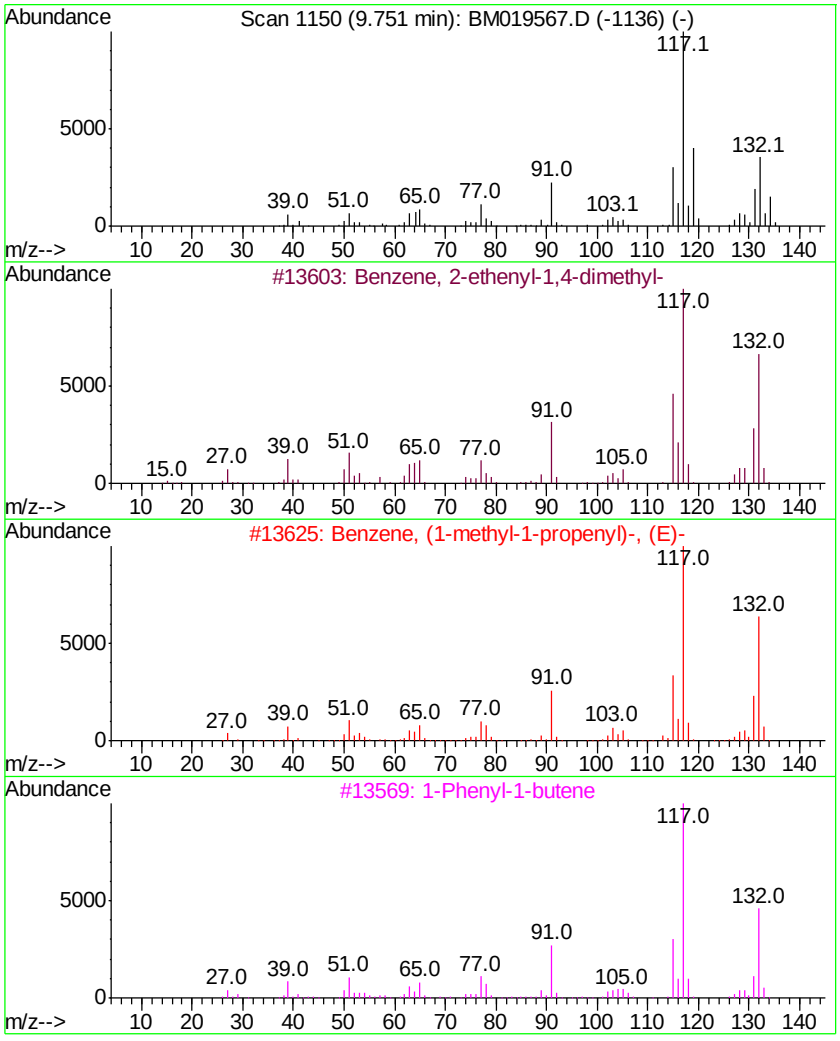
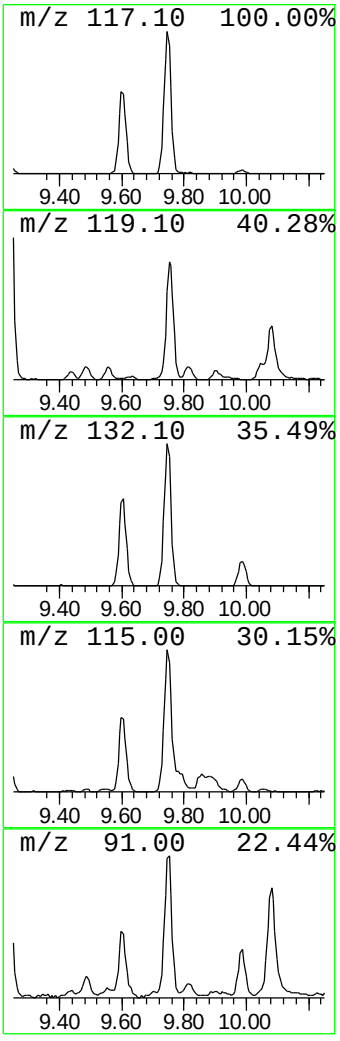
Peak Number 31 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.75	12.09 ng/ul	896781	Naphthalene-d8	10.36

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	95
2	Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	91
3	1-Phenyl-1-butene	132	C10H12	000824-90-8	87
4	Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	83
5	Indan, 1-methyl-	132	C10H12	000767-58-8	81





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM032819\
Data File : BM019567.D
Acq On : 29 Mar 2019 07:04
Operator : JU/SJ
Sample : K2063-16DL 10X
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : DB6S4DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM032219MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

					Internal Standard			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
2-Pentanol, 2-met...	3.38	3.4	ng/ul	181076	1	7.57	1067460	20.0
3-Pentanol, 3-met...	3.63	4.6	ng/ul	245630	1	7.57	1067460	20.0
Cyclopentanol, 1-...	4.12	2.4	ng/ul	127205	1	7.57	1067460	20.0
2-Hexanol, 2-methyl-	4.60	2.7	ng/ul	144162	1	7.57	1067460	20.0
3-Pentanol, 2,3-d...	4.77	3.6	ng/ul	194973	1	7.57	1067460	20.0
1,3-Dimethylcyclo...	4.82	3.3	ng/ul	176165	1	7.57	1067460	20.0
1,3-Cyclopentadie...	5.11	273.2	ng/ul	14581400	1	7.57	1067460	20.0
Benzene, (1-methy...	6.07	11.5	ng/ul	612490	1	7.57	1067460	20.0
N-Benzyl-2-phenet...	6.55	26.4	ng/ul	1407400	1	7.57	1067460	20.0
Benzene, 1-ethyl-...	6.66	113.3	ng/ul	6048850	1	7.57	1067460	20.0
Benzene, 1-ethyl-...	6.72	45.5	ng/ul	2430560	1	7.57	1067460	20.0
Benzene, 1,3,5-tr...	6.79	47.0	ng/ul	2506950	1	7.57	1067460	20.0
Benzene, 1,2,3-tr...	7.22	189.8	ng/ul	10133100	1	7.57	1067460	20.0
Indane	7.93	42.9	ng/ul	2290430	1	7.57	1067460	20.0
Benzene, 1,3-diet...	8.05	3.9	ng/ul	205762	1	7.57	1067460	20.0
unknown-01	8.10	10.1	ng/ul	537667	1	7.57	1067460	20.0
Benzene, 1,4-diet...	8.20	13.2	ng/ul	704470	1	7.57	1067460	20.0
Benzene, 1-methyl...	8.36	2.8	ng/ul	149368	1	7.57	1067460	20.0
Benzene, 2-ethyl-...	8.51	7.7	ng/ul	409808	1	7.57	1067460	20.0
Benzene, 1-methyl...	8.56	6.7	ng/ul	357047	1	7.57	1067460	20.0
unknown-02	8.62	4.5	ng/ul	242908	1	7.57	1067460	20.0
Benzene, 4-ethyl-...	8.99	2.5	ng/ul	183276	2	10.36	1483620	20.0
1,3-Cyclopentadie...	9.18	4.5	ng/ul	332086	2	10.36	1483620	20.0
Benzene, 1,2,3,4-...	9.24	6.3	ng/ul	464460	2	10.36	1483620	20.0
1H-Indene, 2,3-di...	9.60	5.4	ng/ul	402426	2	10.36	1483620	20.0
Benzene, 2-etheny...	9.75	12.1	ng/ul	896781	2	10.36	1483620	20.0