LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM100119\
Data File : BM022875.D
Acq On : 02 Oct 2019 04:21
Operator : JU
Sample : K4932-07
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BFDG9

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-BM092719MA.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	3.305	50	54	69	rVV	2343810	2896229	34.75%	1.890%
2	3.610	102	106	112	rBV	260627	316358	3.80%	0.206%
3	3.728	122	126	131	rBV	229431	285077	3.42%	0.186%
4	3.828	139	143	152	rBV	863601	1102730	13.23%	0.719%
5	4.005	166	173	186	rVB	5184954	7366882	88.39%	4.807%
6	4.269	213	218	224	rBV	469252	602619	7.23%	0.393%
7	4.446	243	248	256	rBV	698657	925091	11.10%	0.604%
8	4.534	256	263	272	rVB3	108274	222843	2.67%	0.145%
9	4.899	318	325	334	rVV	773523	1123023	13.47%	0.733%
10	5.322	391	397	405	rVV	1522871	2188040	26.25%	1.428%
11	5.457	413	420	429	rVV	4219453	8334882	100.00%	5.438%
12	5.763	468	472	476	rVV2	172199	260764	3.13%	0.170%
13	5.822	476	482	494	rVB	3402860	5062842	60.74%	3.303%
14	6.299	558	563	574	rVB	147497	247418	2.97%	0.161%
15	6.787	637	646	653	rBV	379517	684555	8.21%	0.447%
16	6.898	659	665	670	rBV	1697027	2471949	29.66%	1.613%
17	6.957	670	675	679	rBV2	998423	1729024	20.74%	1.128%
18	6.993	679	681	684	rVV	561057	768325	9.22%	0.501%
19	7.034	684	688	699	rVB	1225885	1829587	21.95%	1.194%
20	7.134	699	705	711	rBV	756507	1207003	14.48%	0.788%
21	7.198	711	716	722	rVB	879096	1288791	15.46%	0.841%
22	7.263	722	727	732	rVB3	175261	255349	3.06%	0.167%
23	7.334	733	739	749	rVB	945174	1533045	18.39%	1.000%
24	7.457	749	760	767	rBV	4233742	6822716	81.86%	4.452%
25	7.804	810	819	823	rBV	1095455	1843175	22.11%	1.203%
26	7.869	823	830	834	rBV2	912083	1351968	16.22%	0.882%
27	7.922	834	839	855	rVB	1827140	3314141	39.76%	2.162%
28	8.093	863	868	873	rBV3	115926	227153	2.73%	0.148%
29	8.175	873	882	889	rBV2	967272	1926131	23.11%	1.257%
30	8.240	889	893	898	rVB	322469	480648	5.77%	0.314%
31	8.298	898	903	906	rBV2	191111	293149	3.52%	0.191%
32	8.351	906	912	921	rVB2	360272	714569	8.57%	0.466%
33	8.445	921	928	933	rBV	670739	1049789	12.60%	0.685%
34	8.504	933	938	943	rBV	781970	1221859	14.66%	0.797%

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Misc :
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Instrument : BNA_M
ClientSampleId : BFDG9

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

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Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM092719MA.M
Title : SVOA CALIBRATION

35	8.569	943	949	954	rBV	1076845	1813939	21.76%	1.184%
36	8.610	954	956	964	rVB2	175016	266387	3.20%	0.174%
37	8.757	975	981	985	rBV	334990	513368	6.16%	0.335%
38	8.810	985	990	997	rVB	354640	553380	6.64%	0.361%
39	8.910	997	1007	1011	rBV	767923	1414464	16.97%	0.923%
40	8.957	1011	1015	1018	rBV	738267	1091014	13.09%	0.712%
41	8.987	1018	1020	1026	rVB2	432757	580061	6.96%	0.378%
42	9.045	1026	1030	1034	rVB2	222565	327651	3.93%	0.214%
43	9.240	1057	1063	1070	rVB2	210974	369903	4.44%	0.241%
44	9.428	1090	1095	1100	rVB	393535	602559	7.23%	0.393%
45	9.492	1100	1106	1115	rBV	624645	1142272	13.70%	0.745%
46	9.675	1129	1137	1144	rBV	818559	1448517	17.38%	0.945%
47	9.769	1144	1153	1157	rBV	908674	1656003	19.87%	1.080%
48	9.804	1157	1159	1163	rVB2	398155	424266	5.09%	0.277%
49	9.845	1163	1166	1175	rVB	485837	735893	8.83%	0.480%
50	9.928	1175	1180	1183	rBV	108566	179416	2.15%	0.117%
51	10.004	1185	1193	1202	rBV5	1330407	3854933	46.25%	2.515%
52	10.075	1202	1205	1214	rVB	799733	1234867	14.82%	0.806%
53	10.216	1222	1229	1238	rVB2	1464963	2934907	35.21%	1.915%
54	10.298	1238	1243	1251	rVB2	174492	298735	3.58%	0.195%
55	10.375	1251	1256	1264	rBV	230224	397920	4.77%	0.260%
56	10.463	1265	1271	1276	rBV	376589	680351	8.16%	0.444%
57	10.592	1284	1293	1297	rBV	1594202	2916973	35.00%	1.903%
58	10.645	1297	1302	1309	rVB	4675183	7893696	94.71%	5.150%
59	10.722	1309	1315	1320	rBV	877869	1772718	21.27%	1.157%
60	10.945	1347	1353	1362	rVB	452757	867005	10.40%	0.566%
61	11.128	1378	1384	1390	rBV	196983	337271	4.05%	0.220%
62	11.204	1390	1397	1402	rBV3	119934	249787	3.00%	0.163%
63	11.422	1421	1434	1440	rBV4	254895	700828	8.41%	0.457%
64	11.510	1444	1449	1454	rVB	185880	301796	3.62%	0.197%
65	11.616	1462	1467	1473	rBV2	145048	280123	3.36%	0.183%
66	11.704	1473	1482	1490	rVB2	242735	535693	6.43%	0.350%
67	11.933	1509	1521	1526	rBV2	199055	439963	5.28%	0.287%
68	12.151	1553	1558	1563	rBV2	146230	260262	3.12%	0.170%
69	12.257	1569	1576	1589	rVB	2500068	3919986	47.03%	2.558%
70	12.375	1589	1596	1600	rBV3	87295	214115	2.57%	0.140%
71	12.475	1605	1613	1619	rVB	1316514	2200333	26.40%	1.436%

LSC Area Percent Report

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Instrument : BNA_M
ClientSampleId : BFDG9

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

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Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM092719MA.M
Title : SVOA CALIBRATION

72	12.722	1647	1655	1660	rVB4	116691	264639	3.18%	0.173%
73	13.022	1702	1706	1717	rVB	165578	302770	3.63%	0.198%
74	13.280	1744	1750	1758	rVV2	361603	640413	7.68%	0.418%
75	13.351	1758	1762	1767	rVV	281651	391452	4.70%	0.255%
76	13.463	1775	1781	1794	rVB4	159235	544028	6.53%	0.355%
77	13.616	1800	1807	1812	rVB4	179266	432776	5.19%	0.282%
78	13.757	1826	1831	1836	rVV	240403	429930	5.16%	0.281%
79	13.816	1836	1841	1842	rVV	202686	298106	3.58%	0.195%
80	13.845	1842	1846	1850	rVV	2005004	2908729	34.90%	1.898%
81	13.892	1850	1854	1865	rVB	1086182	1600206	19.20%	1.044%
82	13.992	1865	1871	1875	rBV2	136147	249086	2.99%	0.163%
83	14.127	1888	1894	1905	rVB	2404620	3839754	46.07%	2.505%
84	14.439	1940	1947	1954	rVV	2252877	3392859	40.71%	2.214%
85	14.633	1976	1980	1986	rVB	209015	298017	3.58%	0.194%
86	15.427	2109	2115	2121	rBV	2994194	4425018	53.09%	2.887%
87	15.539	2129	2134	2142	rVB	886488	1266648	15.20%	0.826%
88	16.427	2281	2285	2288	rBV	178219	226533	2.72%	0.148%
89	16.827	2349	2353	2359	rVB3	163860	232444	2.79%	0.152%
90	17.174	2406	2412	2417	rBV2	2251601	3202513	38.42%	2.090%
91	17.274	2423	2429	2435	rBV	2924546	4159340	49.90%	2.714%
92	17.574	2477	2480	2489	rVB	126793	188870	2.27%	0.123%
93	18.074	2561	2565	2572	rBV	535066	759218	9.11%	0.495%
94	19.357	2779	2783	2793	rVB	253504	372923	4.47%	0.243%
95	19.568	2813	2819	2828	rBV	3084823	4374303	52.48%	2.854%
96	21.045	3067	3070	3071	rBV	215812	229072	2.75%	0.149%
97	21.068	3071	3074	3078	rVB	426204	514407	6.17%	0.336%
98	21.356	3118	3123	3130	rBV	2260533	2931359	35.17%	1.913%
99	23.474	3477	3483	3498	rVB	2153386	4403239	52.83%	2.873%
100	23.615	3501	3507	3515	rVB	1567875	3028049	36.33%	1.976%

Sum of corrected areas: 153265780

LSC Report - Integrated Chromatogram

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM100119\
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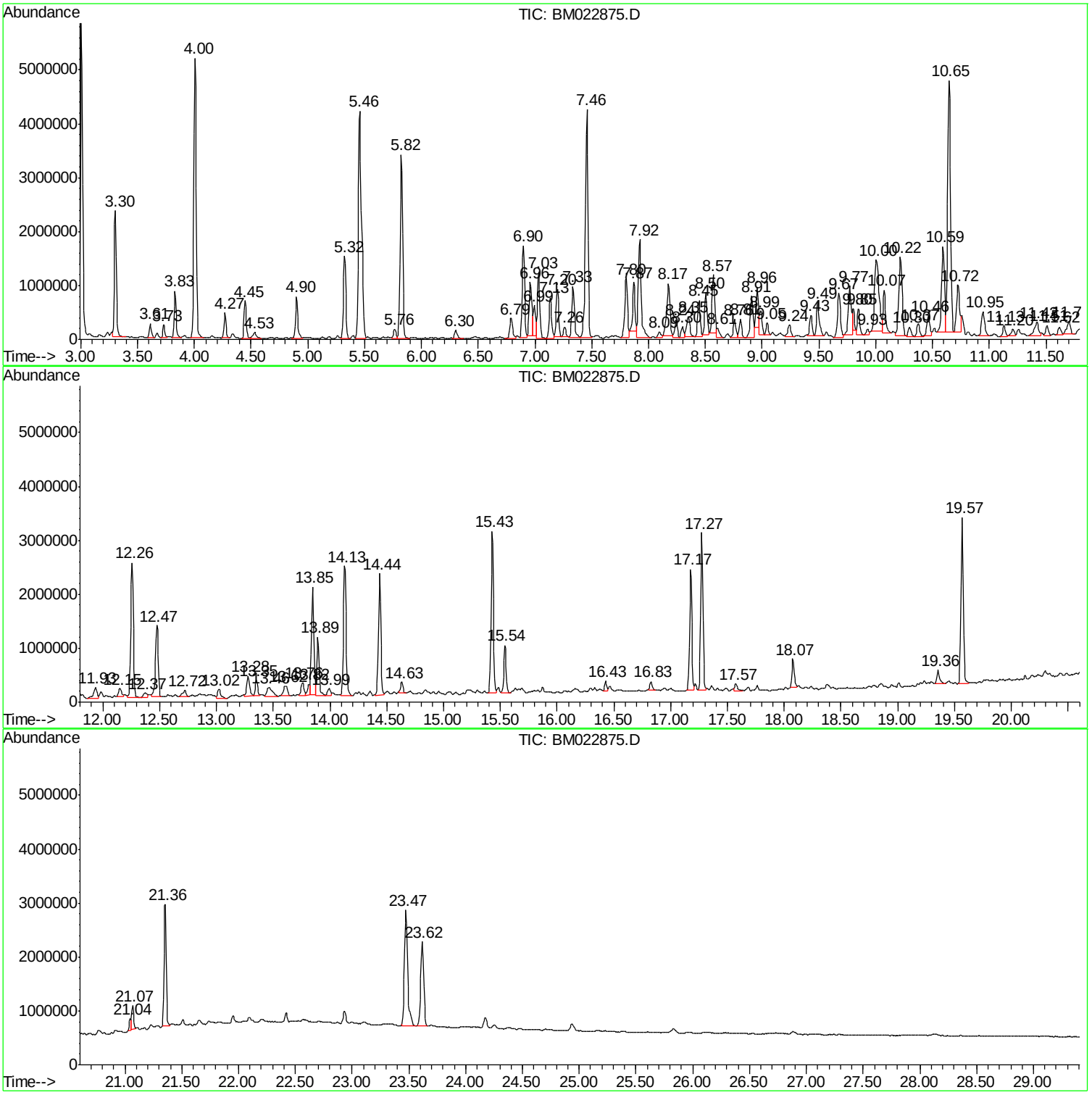
ALS Vial : 21 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BFDG9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM092719MA.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM100119\
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Quant Title : SVOA CALIBRATION

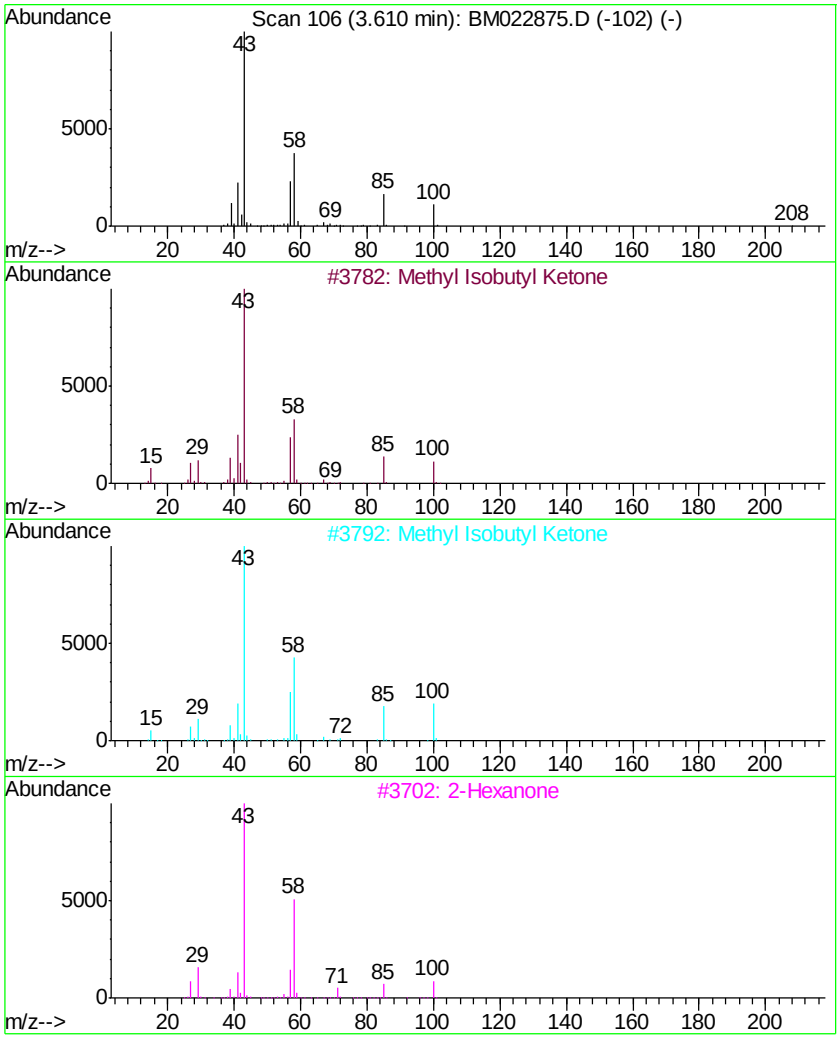
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

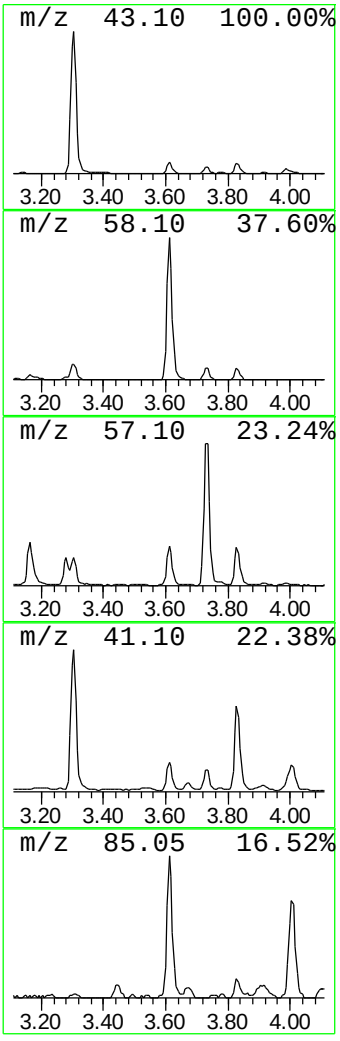
****

Peak Number 1 Methyl Isobutyl Ketone Concentration Rank 33

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.61	3.43 ng/ul	316358	1,4-Dichlorobenzene-d4	7.80

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Methyl Isobutyl Ketone	100	C6H12O	000108-10-1	80
2	Methyl Isobutyl Ketone	100	C6H12O	000108-10-1	72
3	2-Hexanone	100	C6H12O	000591-78-6	72
4	Methyl Isobutyl Ketone	100	C6H12O	000108-10-1	64
5	2-Hexanone	100	C6H12O	000591-78-6	52





Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM100119\
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM092719MA.M
Quant Title : SVOA CALIBRATION

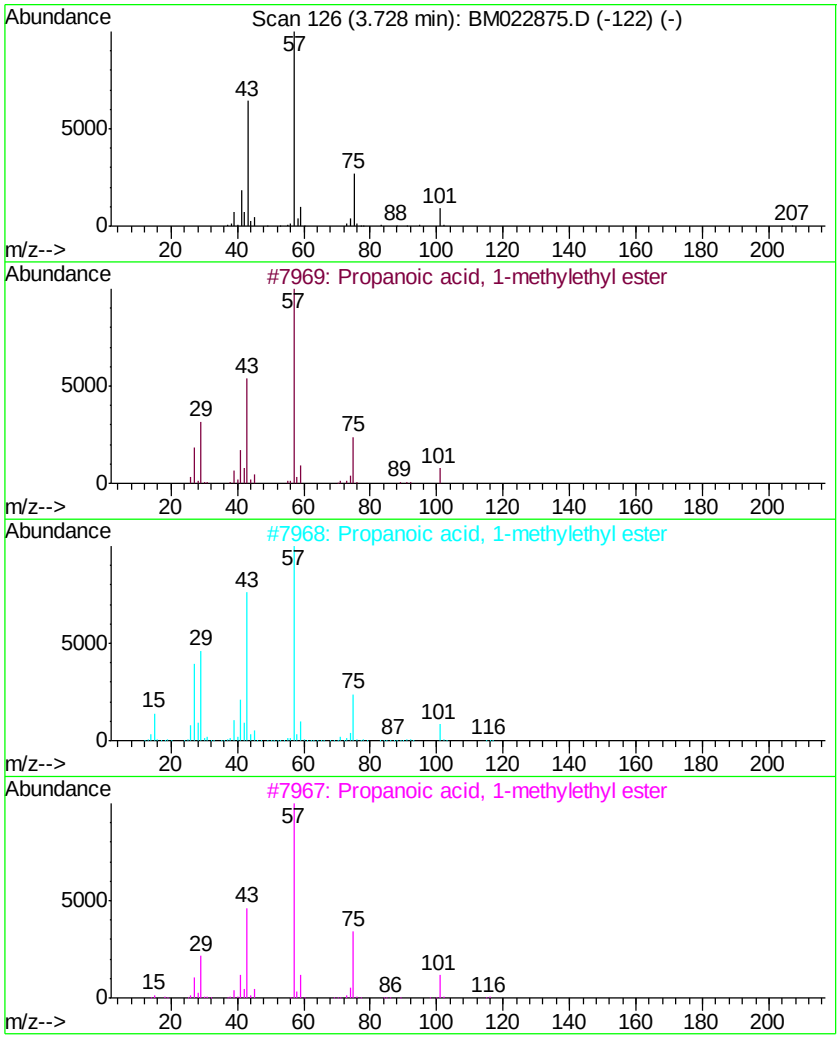
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

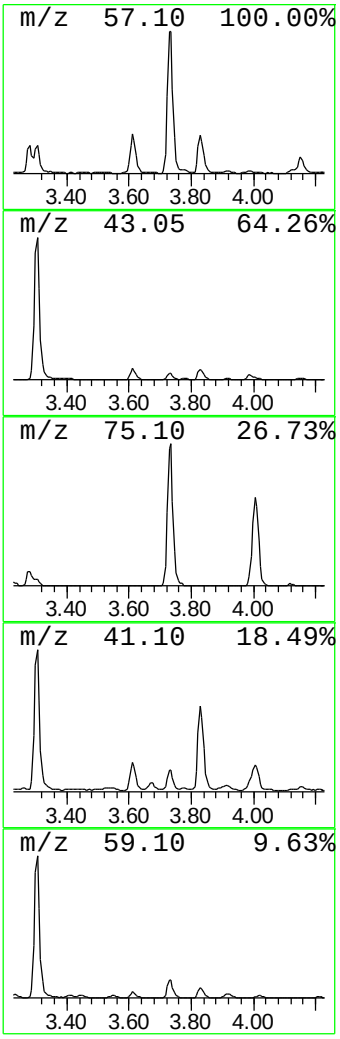
****

Peak Number 2 Propanoic acid, 1-methyleth... Concentration Rank 36

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.73	3.09 ng/ul	285077	1,4-Dichlorobenzene-d4	7.80

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Propanoic acid, 1-methylethyl ester	116	C6H12O2	000637-78-5	83
2	Propanoic acid, 1-methylethyl ester	116	C6H12O2	000637-78-5	83
3	Propanoic acid, 1-methylethyl ester	116	C6H12O2	000637-78-5	64
4	Propanoic acid, 1-methylethyl ester	116	C6H12O2	000637-78-5	42
5	Propane, 1-(1-methylethoxy)-	102	C6H14O	000627-08-7	23





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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM092719MA.M
Quant Title : SVOA CALIBRATION

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

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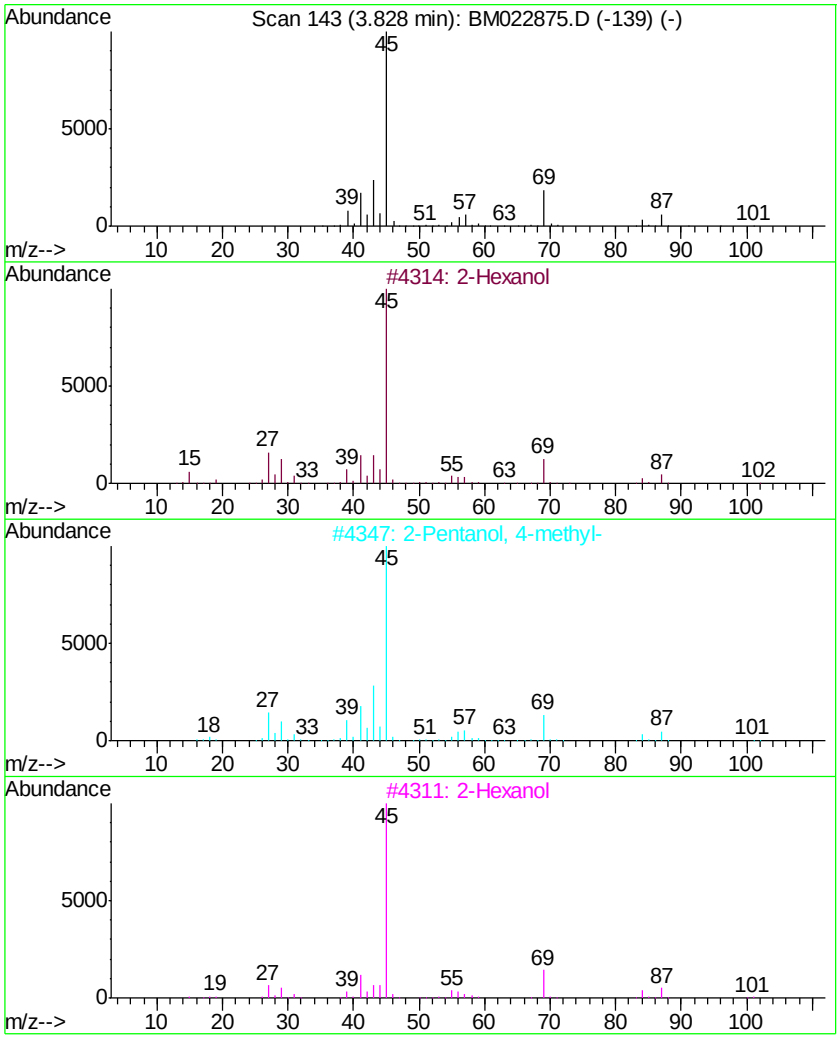
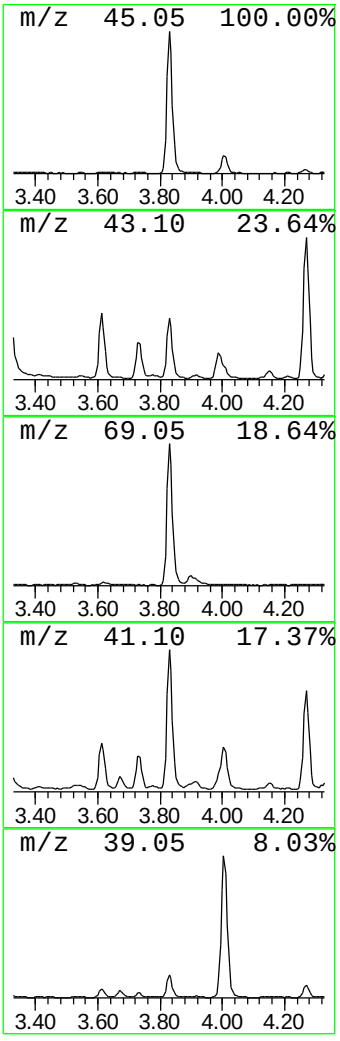
Peak Number 3 2-Hexanol Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.83	11.97 ng/ul	1102730	1,4-Dichlorobenzene-d4	7.80

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Hexanol	102	C6H14O	000626-93-7	83
2	2-Pentanol, 4-methyl-	102	C6H14O	000108-11-2	83
3	2-Hexanol	102	C6H14O	000626-93-7	78
4	2-Hexanol, (R)-	102	C6H14O	026549-24-6	78
5	2-Hexanol	102	C6H14O	000626-93-7	64





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Instrument : BNA_M
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM092719MA.M
Quant Title : SVOA CALIBRATION

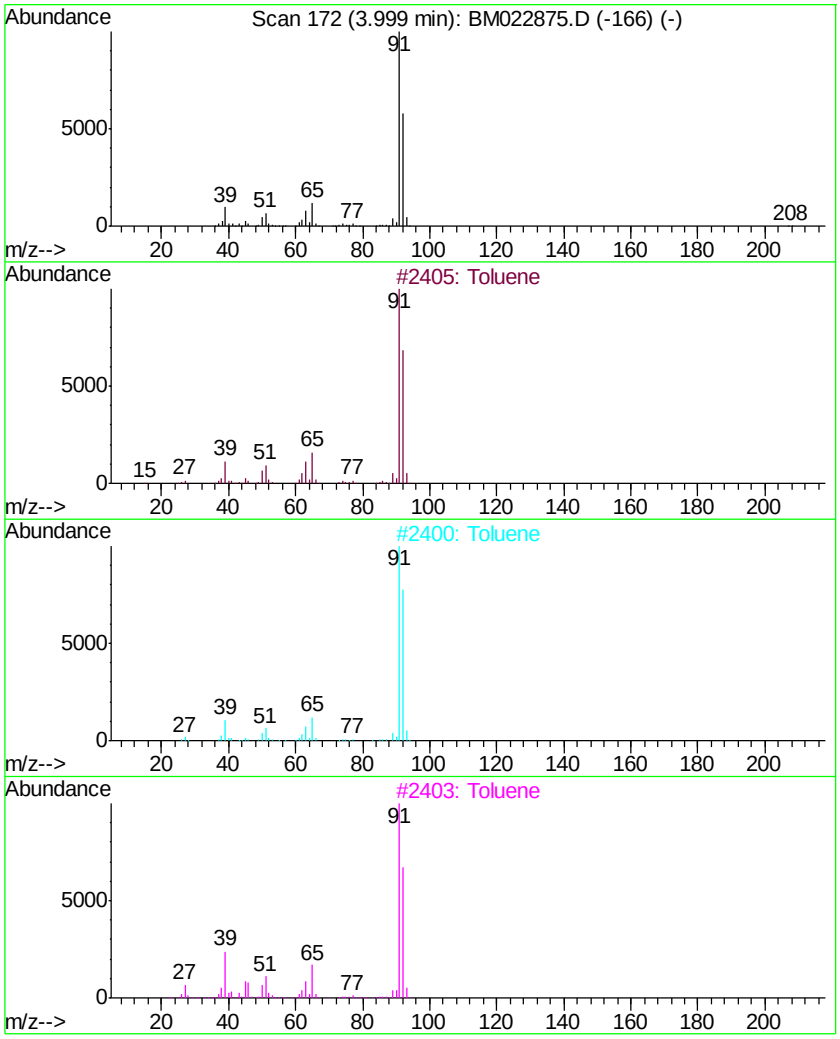
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

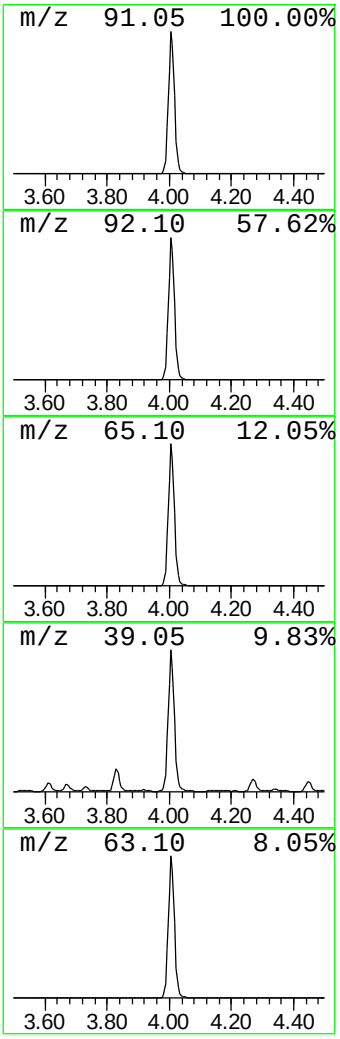
****

Peak Number 4 Toluene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.00	79.94 ng/ul	7366880	1,4-Dichlorobenzene-d4	7.80

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Toluene	92	C7H8	000108-88-3	95
2	Toluene	92	C7H8	000108-88-3	95
3	Toluene	92	C7H8	000108-88-3	91
4	Toluene	92	C7H8	000108-88-3	91
5	1,3,5-Cycloheptatriene	92	C7H8	000544-25-2	87





Library Search Compound Report

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Acq On : 02 Oct 2019 04:21
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Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BFDG9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM092719MA.M
Quant Title : SVOA CALIBRATION

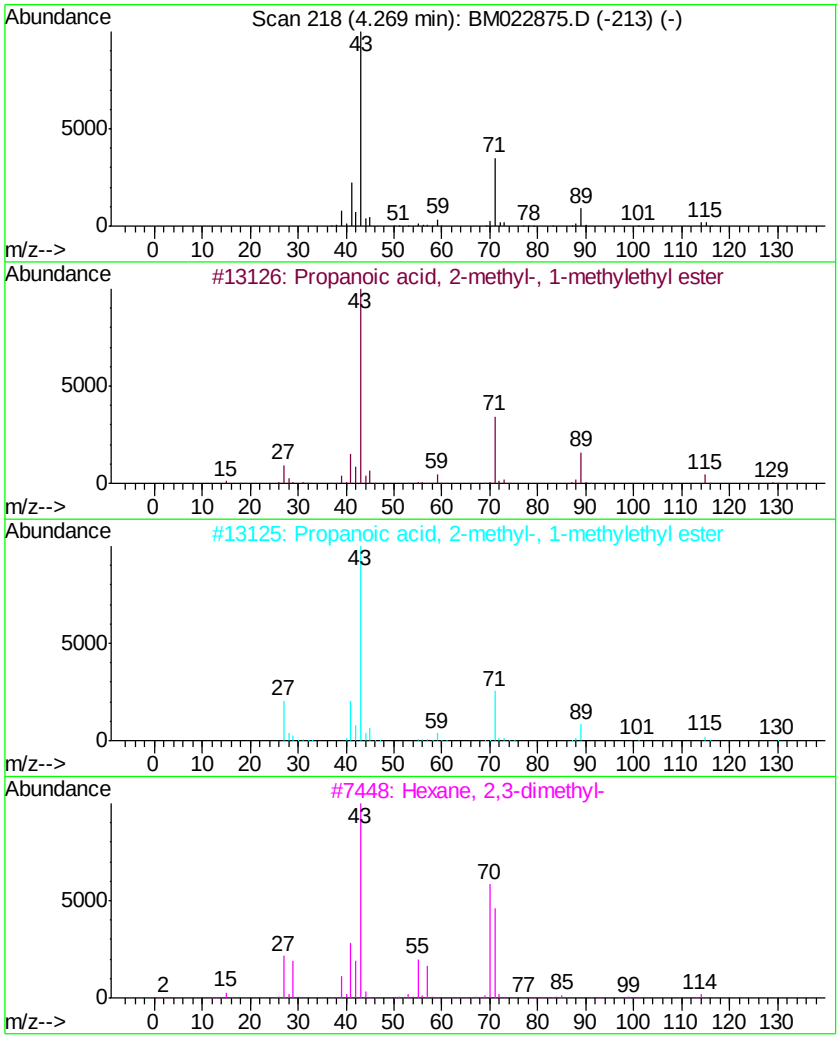
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

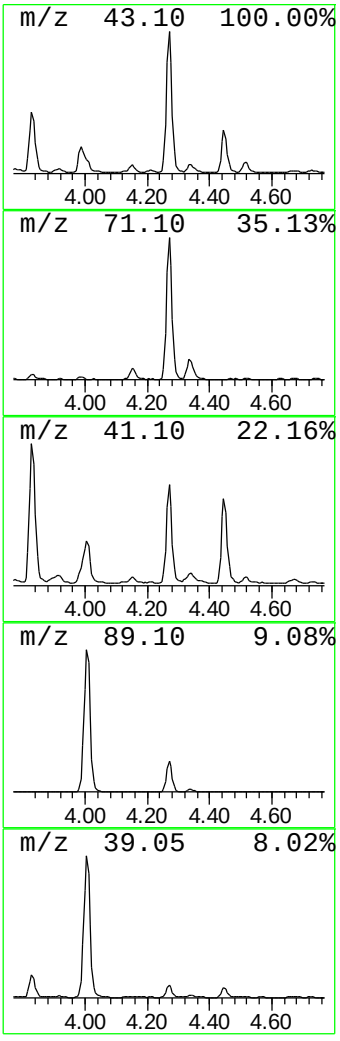
****

Peak Number 5 Propanoic acid, 2-methyl-, ... Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.27	6.54 ng/ul	602619	1,4-Dichlorobenzene-d4	7.80

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Propanoic acid, 2-methyl-, 1-met...	130	C7H14O2	000617-50-5	83
2	Propanoic acid, 2-methyl-, 1-met...	130	C7H14O2	000617-50-5	64
3	Hexane, 2,3-dimethyl-	114	C8H18	000584-94-1	50
4	Propanoic acid, 2-methyl-, 3-met...	158	C9H18O2	002050-01-3	42
5	Propanoic acid, 2-methyl-, 1-met...	130	C7H14O2	000617-50-5	40





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Acq On : 02 Oct 2019 04:21
Operator : JU
Sample : K4932-07
Misc :
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Instrument : BNA_M
ClientSampleId : BFDG9

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Quant Title : SVOA CALIBRATION

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

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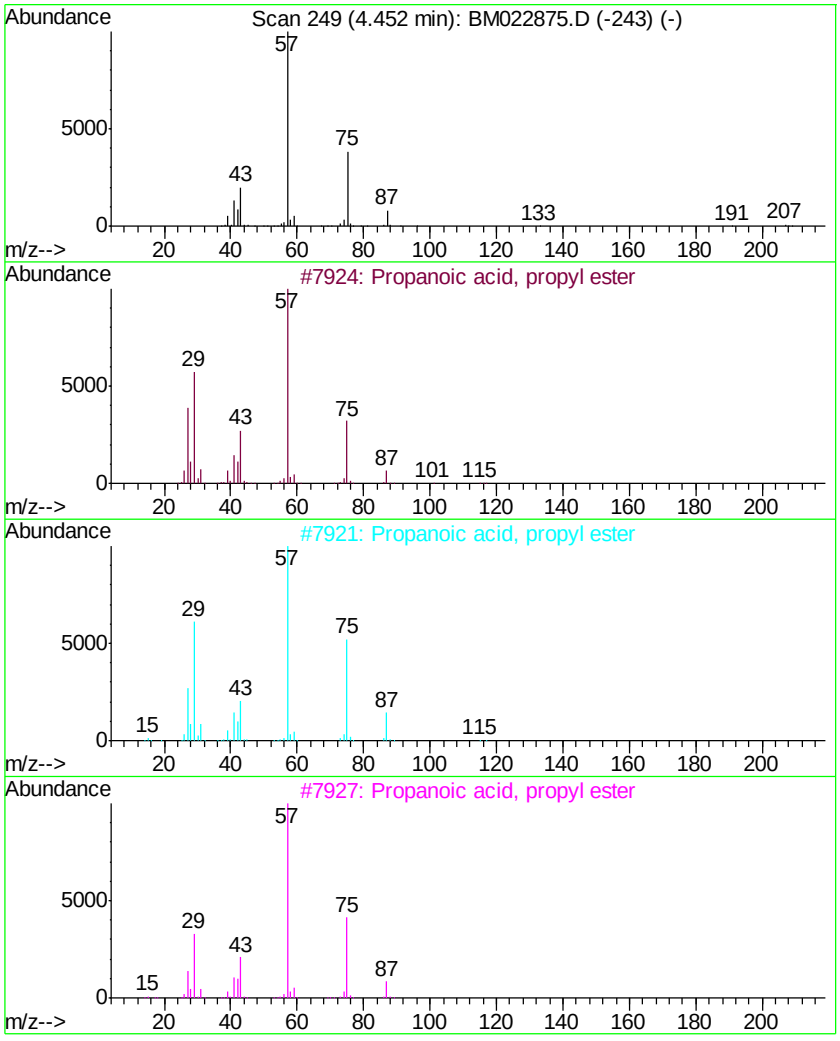
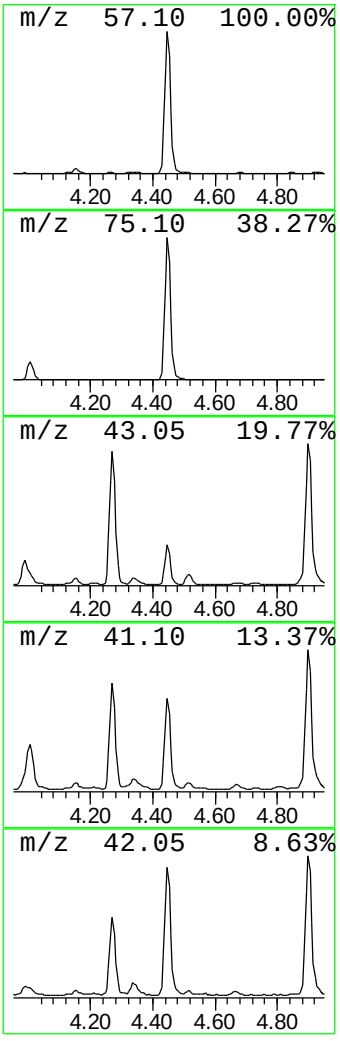
Peak Number 6 Propanoic acid, propyl ester Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.45	10.04 ng/ul	925091	1,4-Dichlorobenzene-d4	7.80

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Propanoic acid, propyl ester	116	C6H12O2	000106-36-5	83
2	Propanoic acid, propyl ester	116	C6H12O2	000106-36-5	83
3	Propanoic acid, propyl ester	116	C6H12O2	000106-36-5	78
4	Propanoic acid, propyl ester	116	C6H12O2	000106-36-5	64
5	Propanoic acid, butyl ester	130	C7H14O2	000590-01-2	39





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Instrument : BNA_M
ClientSampleId : BFDG9

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Quant Title : SVOA CALIBRATION

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

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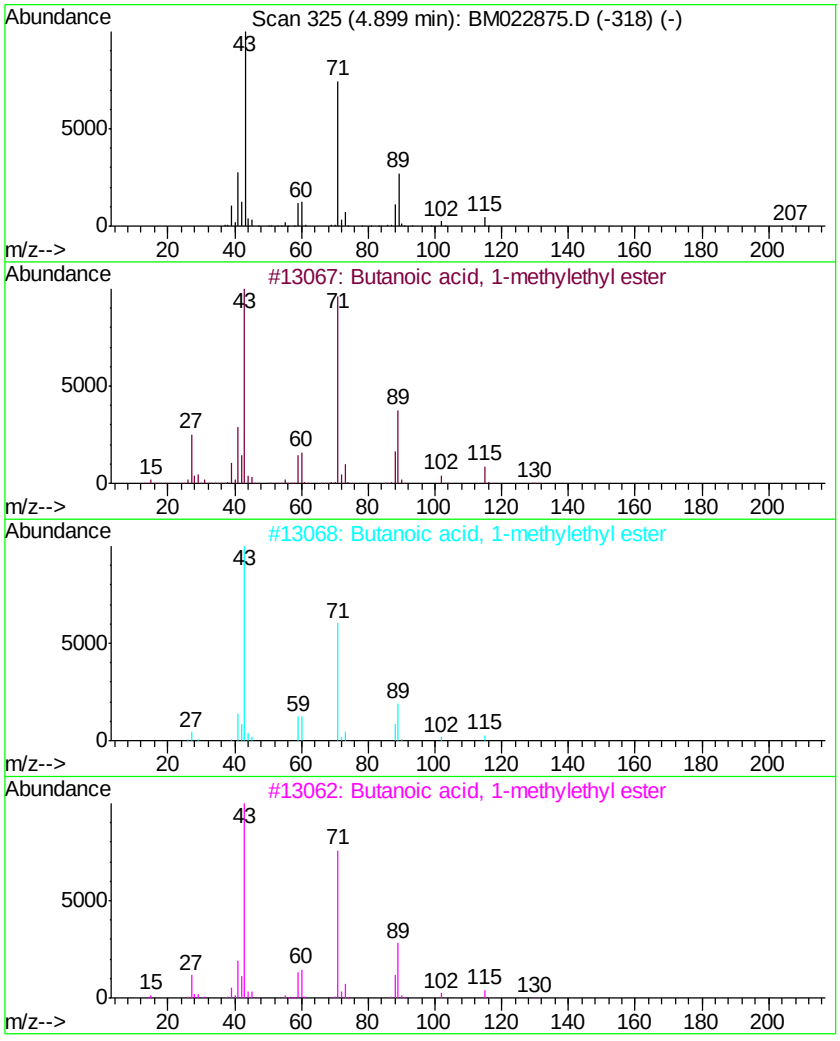
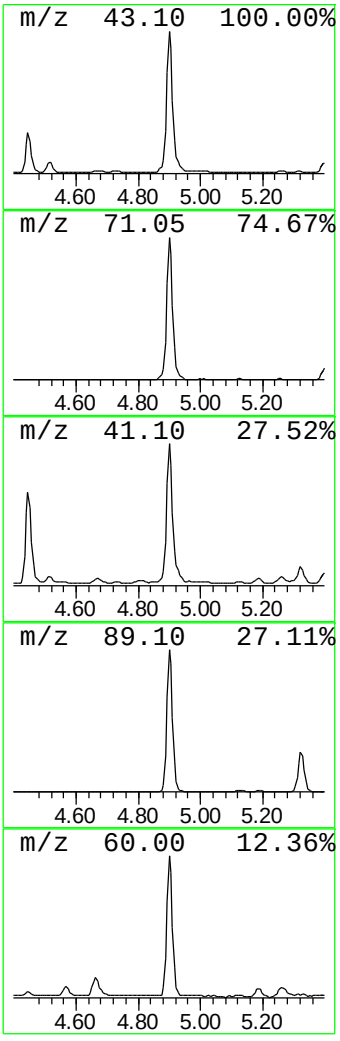
Peak Number 8 Butanoic acid, 1-methylethy... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.90	12.19 ng/ul	1123020	1,4-Dichlorobenzene-d4	7.80

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butanoic acid, 1-methylethyl ester	130	C7H14O2	000638-11-9	90
2	Butanoic acid, 1-methylethyl ester	130	C7H14O2	000638-11-9	90
3	Butanoic acid, 1-methylethyl ester	130	C7H14O2	000638-11-9	90
4	Butanoic acid, 1-methylethyl ester	130	C7H14O2	000638-11-9	53
5	Butanoic acid, propyl ester	130	C7H14O2	000105-66-8	50





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ALS Vial : 21 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BFDG9

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Quant Title : SVOA CALIBRATION

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

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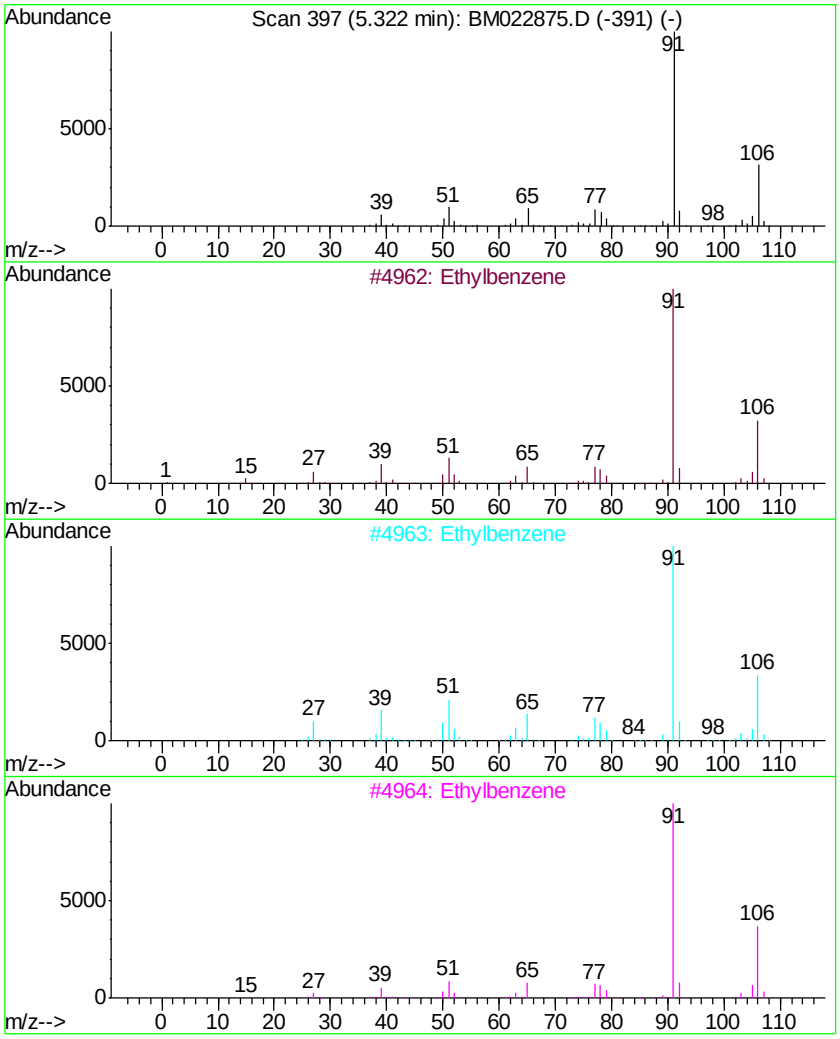
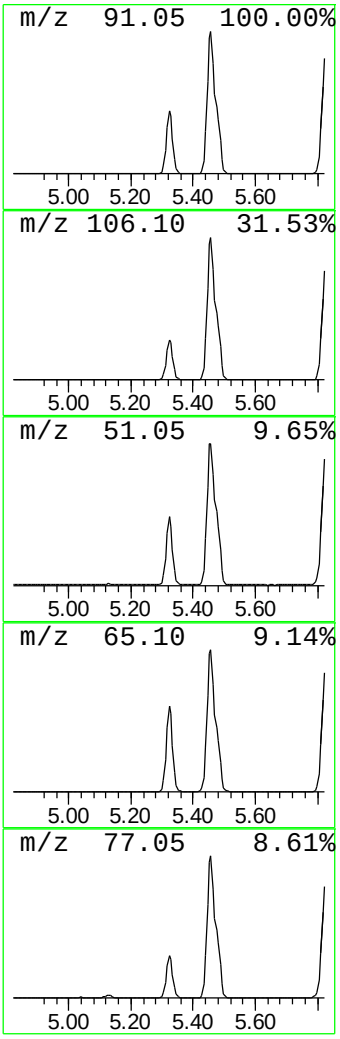
Peak Number 9 Ethylbenzene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.32	23.74 ng/ul	2188040	1,4-Dichlorobenzene-d4	7.80

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethylbenzene	106	C8H10	000100-41-4	94
2	Ethylbenzene	106	C8H10	000100-41-4	91
3	Ethylbenzene	106	C8H10	000100-41-4	90
4	Ethylbenzene	106	C8H10	000100-41-4	83
5	p-Xylene	106	C8H10	000106-42-3	72





Library Search Compound Report

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Quant Title : SVOA CALIBRATION

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

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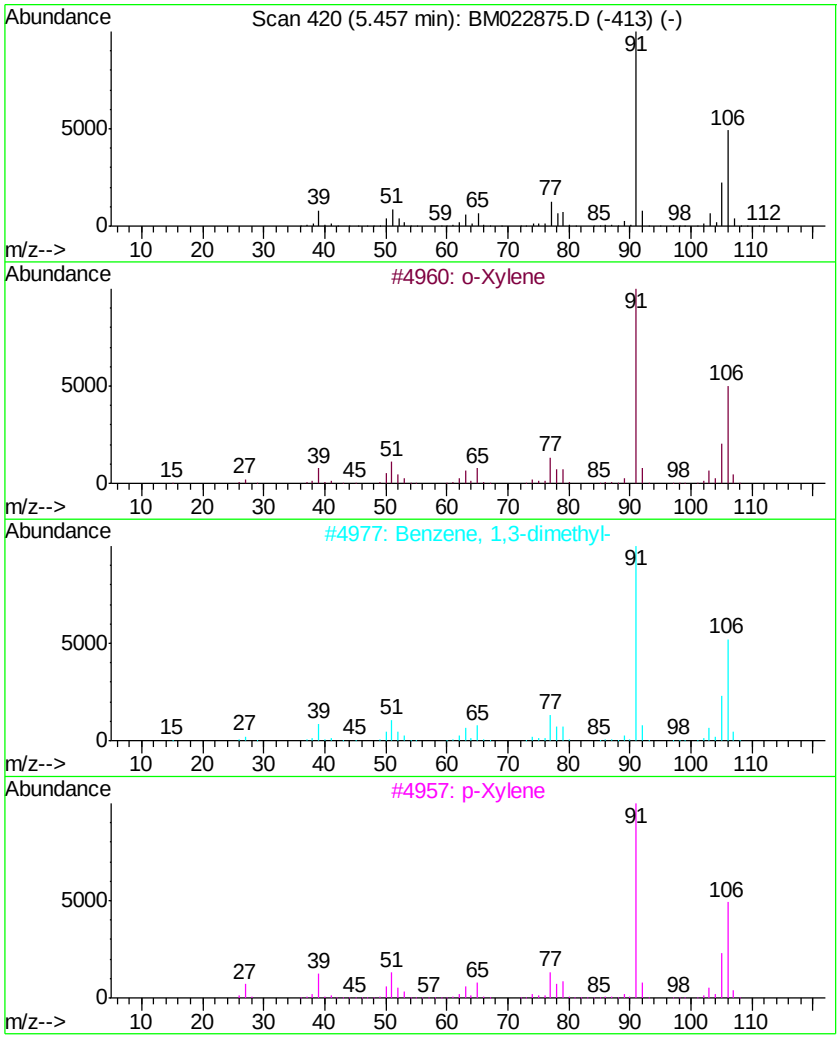
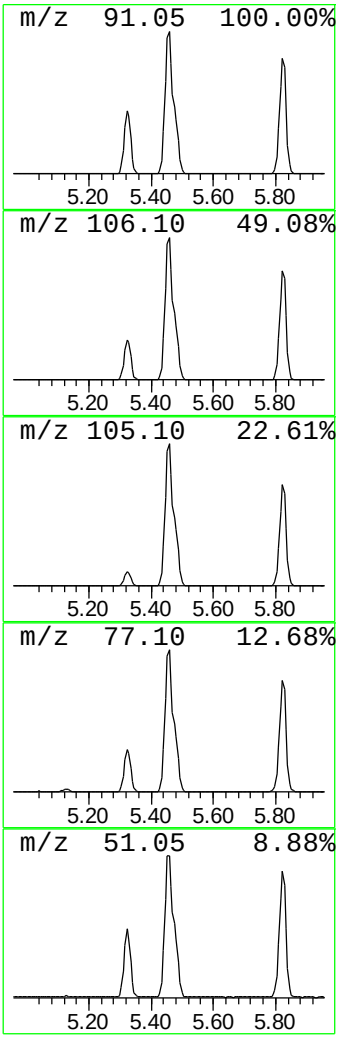
Peak Number 10 o-Xylene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.46	90.44 ng/ul	8334880	1,4-Dichlorobenzene-d4	7.80

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	o-Xylene	106	C8H10	000095-47-6	97
2	Benzene, 1,3-dimethyl-	106	C8H10	000108-38-3	97
3	p-Xylene	106	C8H10	000106-42-3	97
4	p-Xylene	106	C8H10	000106-42-3	97
5	p-Xylene	106	C8H10	000106-42-3	97





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Instrument : BNA_M
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Quant Title : SVOA CALIBRATION

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

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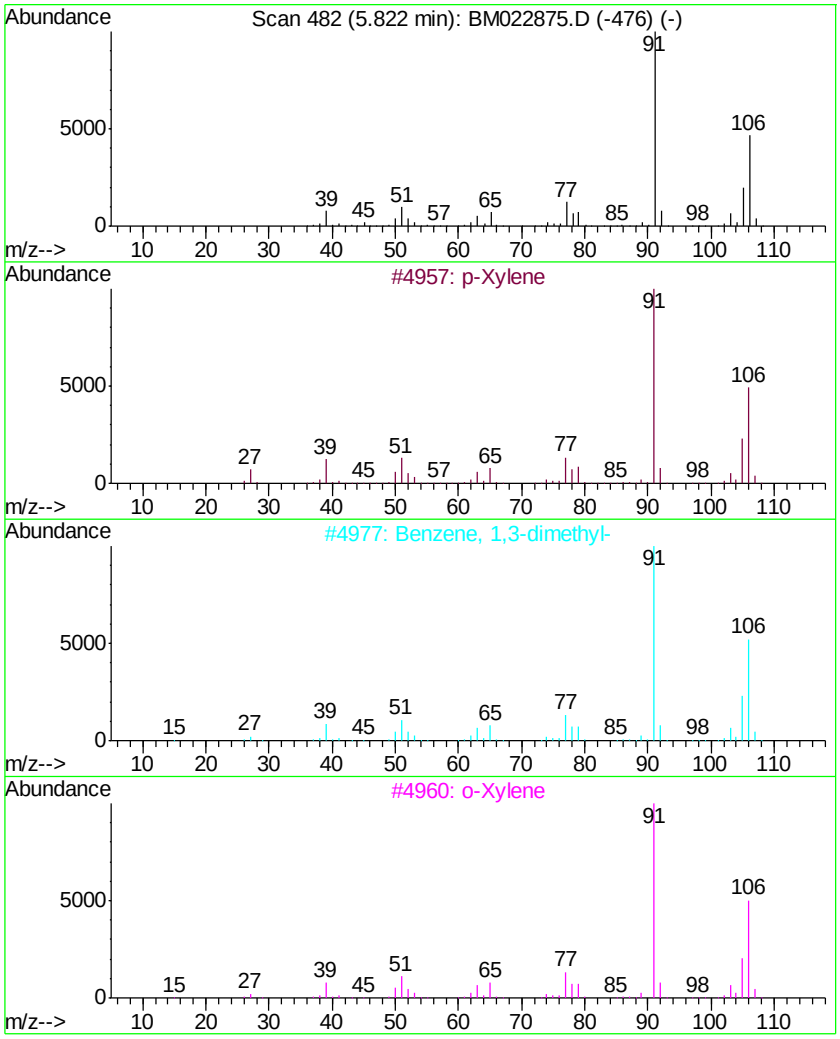
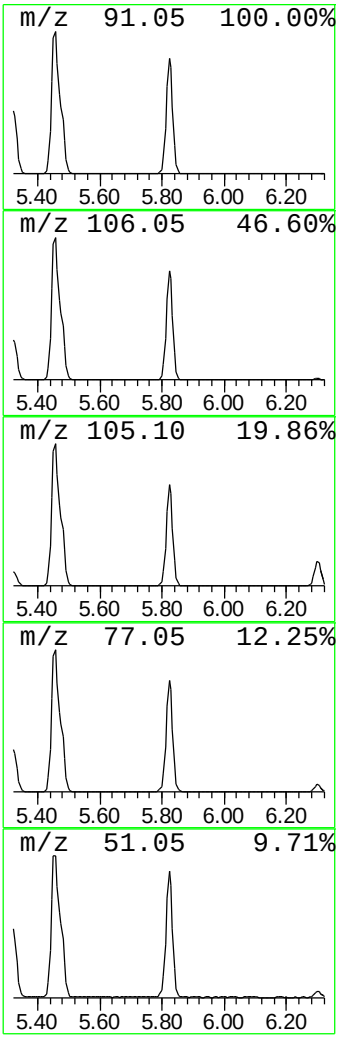
Peak Number 12 p-Xylene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.82	54.94 ng/ul	5062840	1,4-Dichlorobenzene-d4	7.80

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	p-Xylene	106	C8H10	000106-42-3	97
2	Benzene, 1,3-dimethyl-	106	C8H10	000108-38-3	97
3	o-Xylene	106	C8H10	000095-47-6	97
4	p-Xylene	106	C8H10	000106-42-3	95
5	o-Xylene	106	C8H10	000095-47-6	94





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Instrument : BNA_M
ClientSampleId : BFDG9

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Quant Title : SVOA CALIBRATION

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

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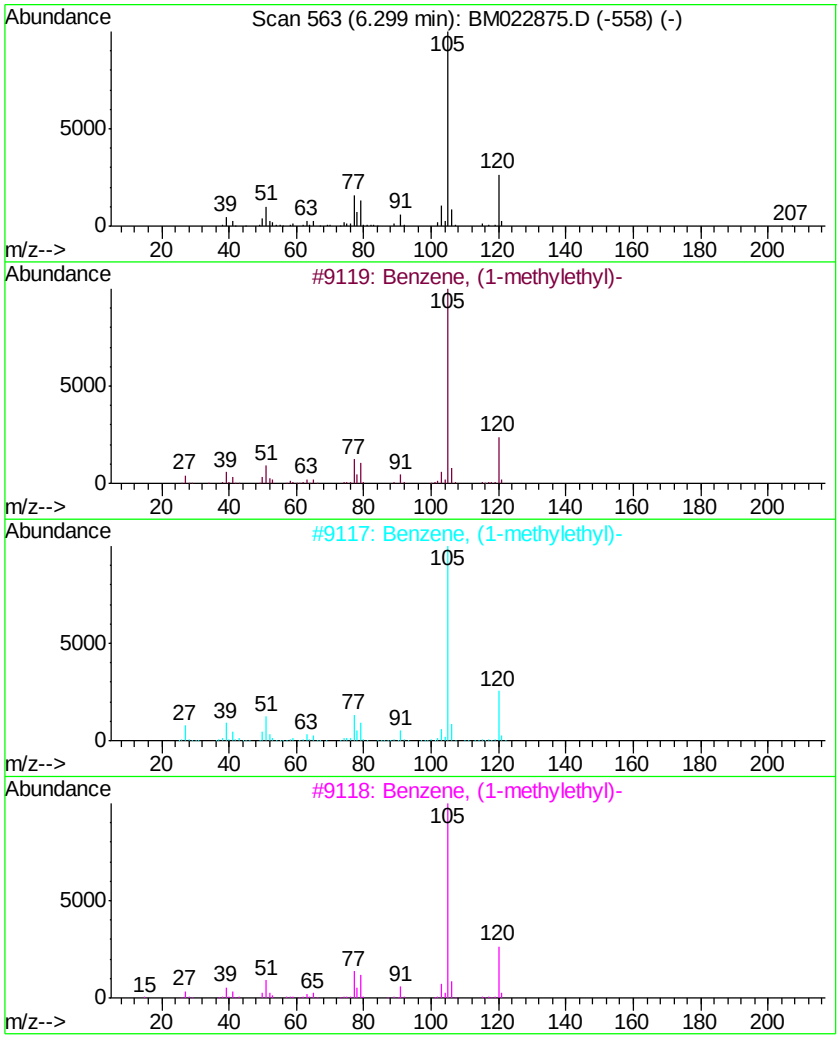
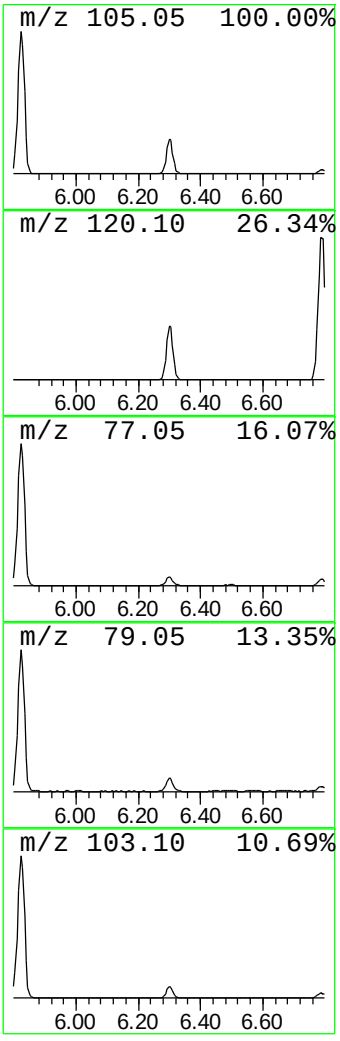
Peak Number 13 Benzene, (1-methylethyl)- Concentration Rank 41

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.30	2.68 ng/ul	247418	1,4-Dichlorobenzene-d4	7.80

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, (1-methylethyl)-	120	C9H12	000098-82-8	94
2	Benzene, (1-methylethyl)-	120	C9H12	000098-82-8	94
3	Benzene, (1-methylethyl)-	120	C9H12	000098-82-8	94
4	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
5	Benzene, (1-methylethyl)-	120	C9H12	000098-82-8	87





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Instrument : BNA_M
ClientSampleId : BFDG9

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Quant Title : SVOA CALIBRATION

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

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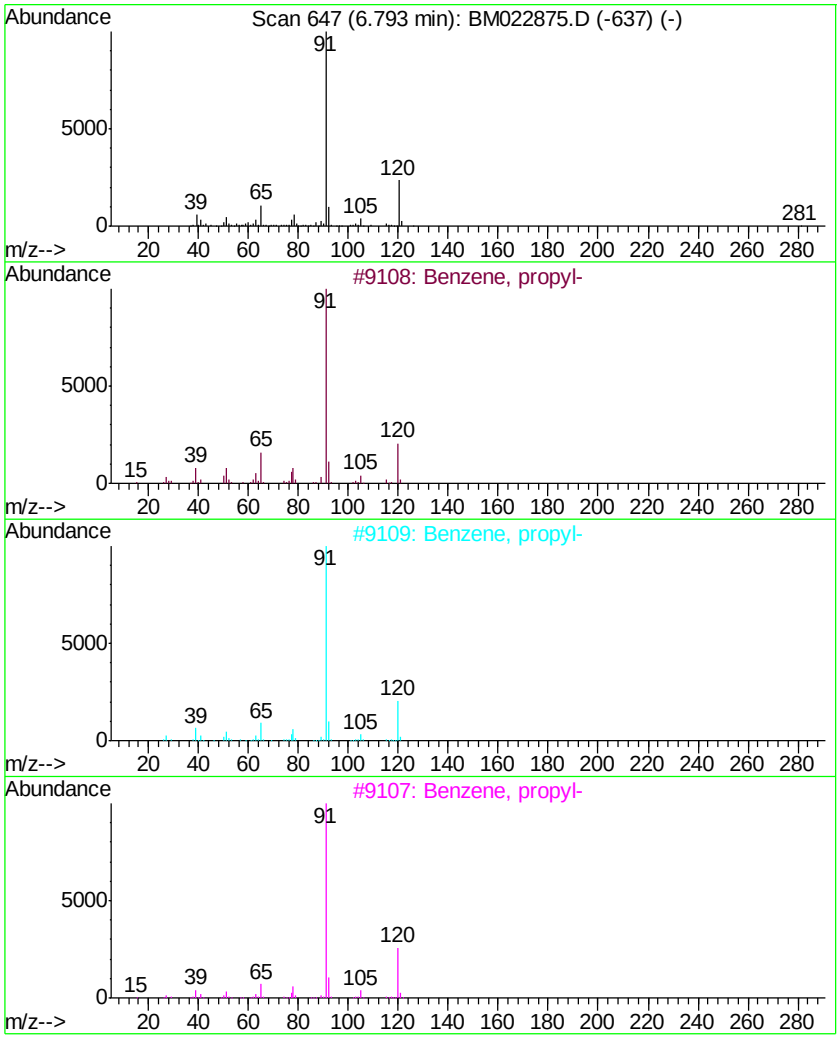
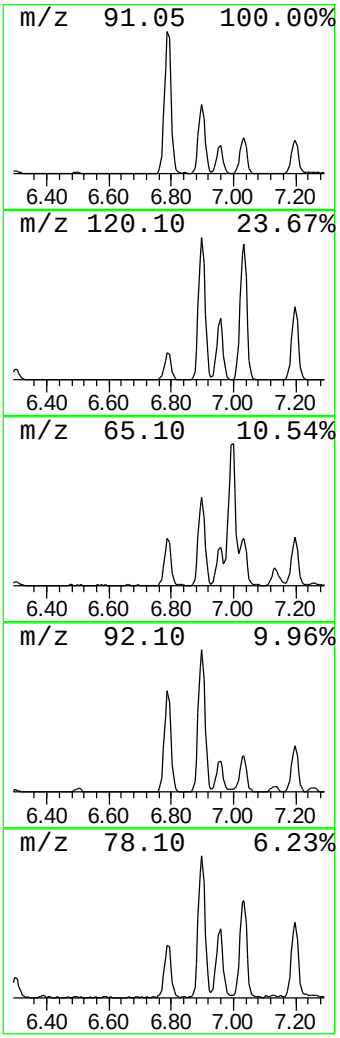
Peak Number 14 Benzene, propyl- Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.79	7.43 ng/ul	684555	1,4-Dichlorobenzene-d4	7.80

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	1,2-Ethanediamine, N,N'-bis(phen...	240	C16H20N2	000140-28-3	72





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Instrument : BNA_M
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Quant Title : SVOA CALIBRATION

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

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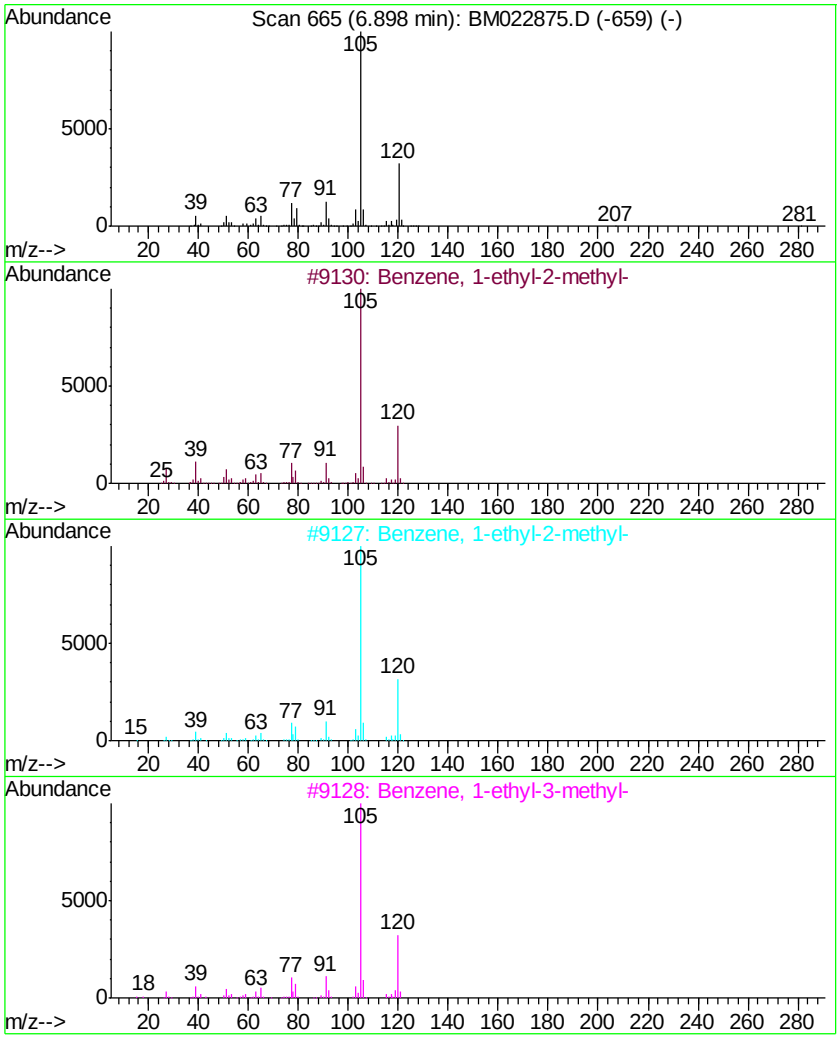
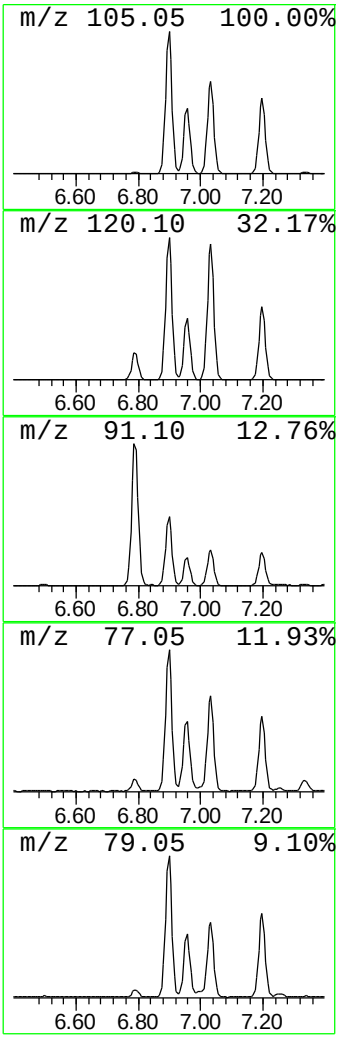
Peak Number 15 Benzene, 1-ethyl-2-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.90	26.82 ng/ul	2471950	1,4-Dichlorobenzene-d4	7.80

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-3-methyl-	120	C9H12	000620-14-4	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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Quant Title : SVOA CALIBRATION

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

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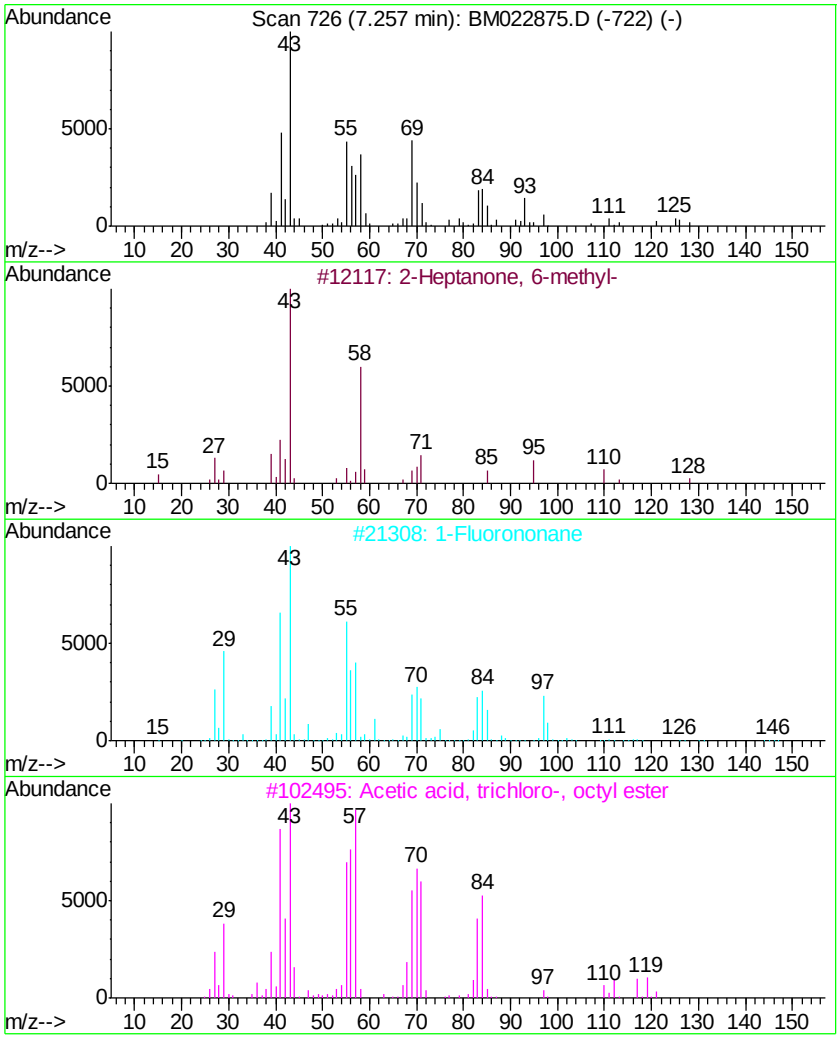
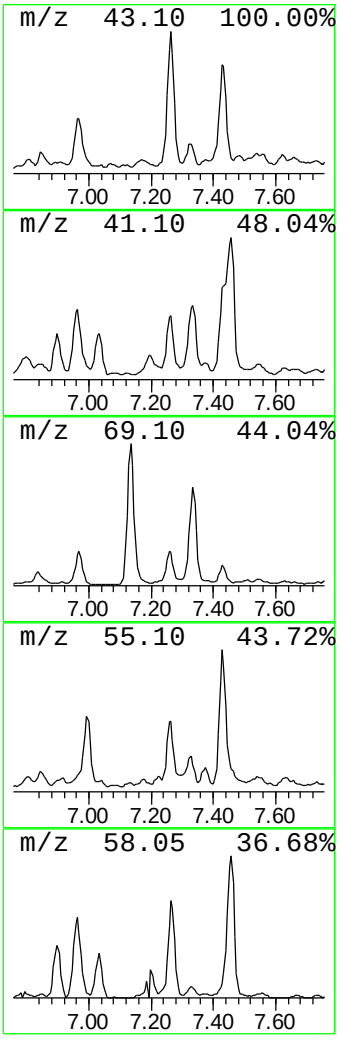
Peak Number 17 unknown-01 Concentration Rank 39

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.26	2.77 ng/ul	255349	1,4-Dichlorobenzene-d4	7.80

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Heptanone, 6-methyl-	128	C8H16O	000928-68-7	43
2	1-Fluorononane	146	C9H19F	000463-18-3	43
3	Acetic acid, trichloro-, octyl e...	274	C10H17Cl3O2	016958-78-4	38
4	1-Fluorononane	146	C9H19F	000463-18-3	37
5	2-Heptanone, 6-methyl-	128	C8H16O	000928-68-7	35





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Quant Title : SVOA CALIBRATION

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

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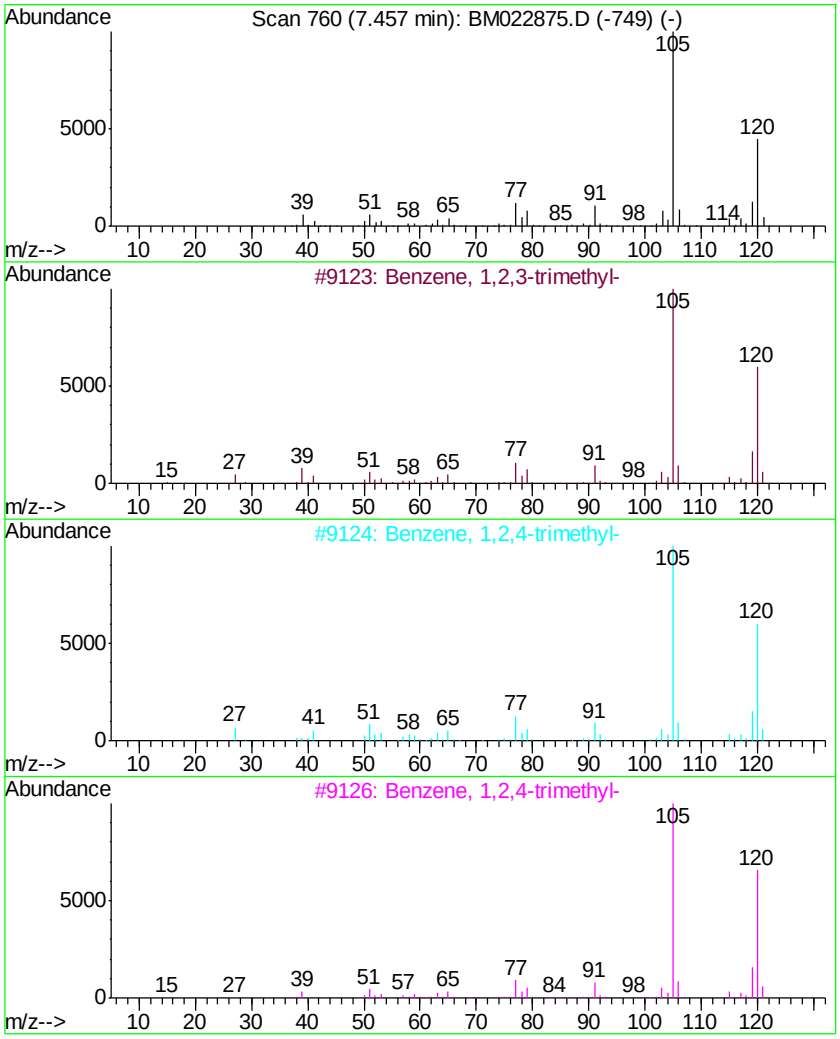
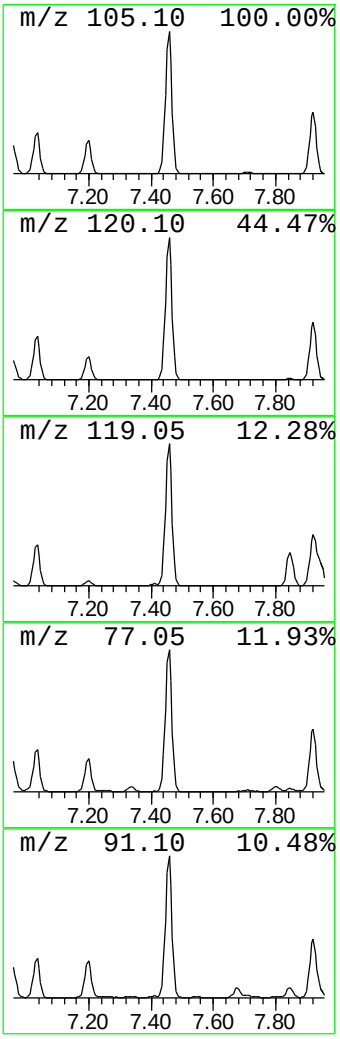
Peak Number 18 Benzene, 1,2,3-trimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.46	74.03 ng/ul	6822720	1,4-Dichlorobenzene-d4	7.80

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	95
3	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94
4	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	94
5	Benzene, 1,3,5-trimethyl-	120	C9H12	000108-67-8	93





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Instrument : BNA_M
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Quant Title : SVOA CALIBRATION

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

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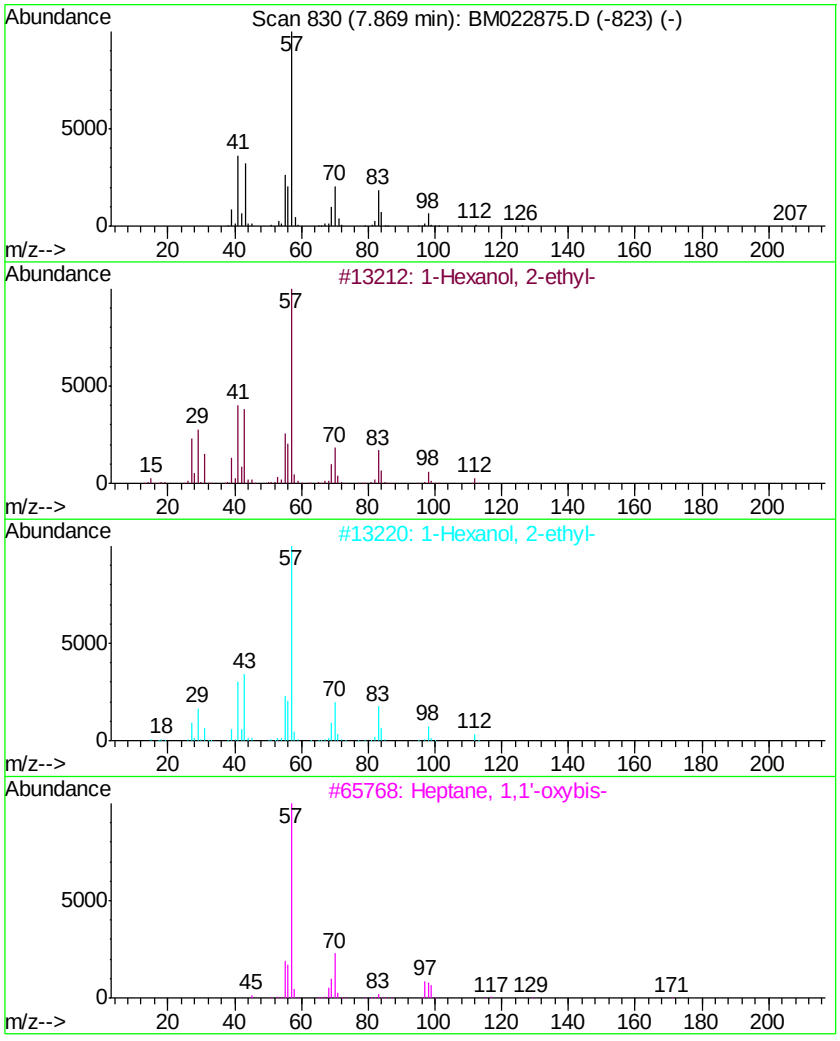
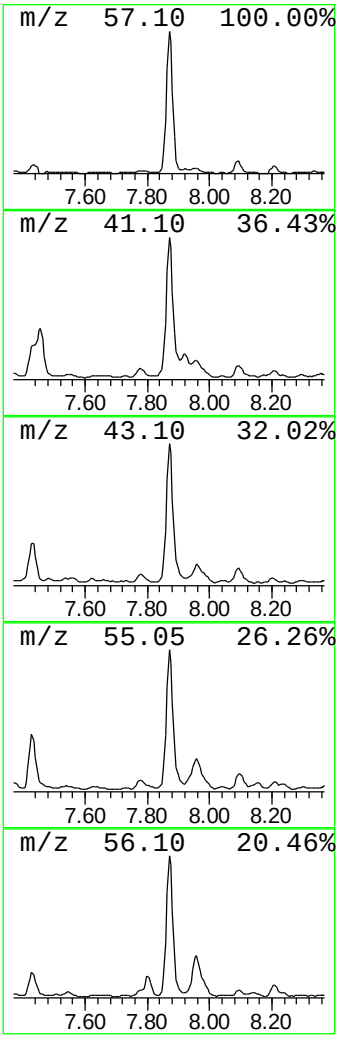
Peak Number 19 1-Hexanol, 2-ethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.87	14.67 ng/ul	1351970	1,4-Dichlorobenzene-d4	7.80

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	83
2	1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	78
3	Heptane, 1,1'-oxybis-	214	C14H30O	000629-64-1	72
4	1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	64
5	Heptane, 1,1'-oxybis-	214	C14H30O	000629-64-1	56





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ALS Vial : 21 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BFDG9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM092719MA.M
Quant Title : SVOA CALIBRATION

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

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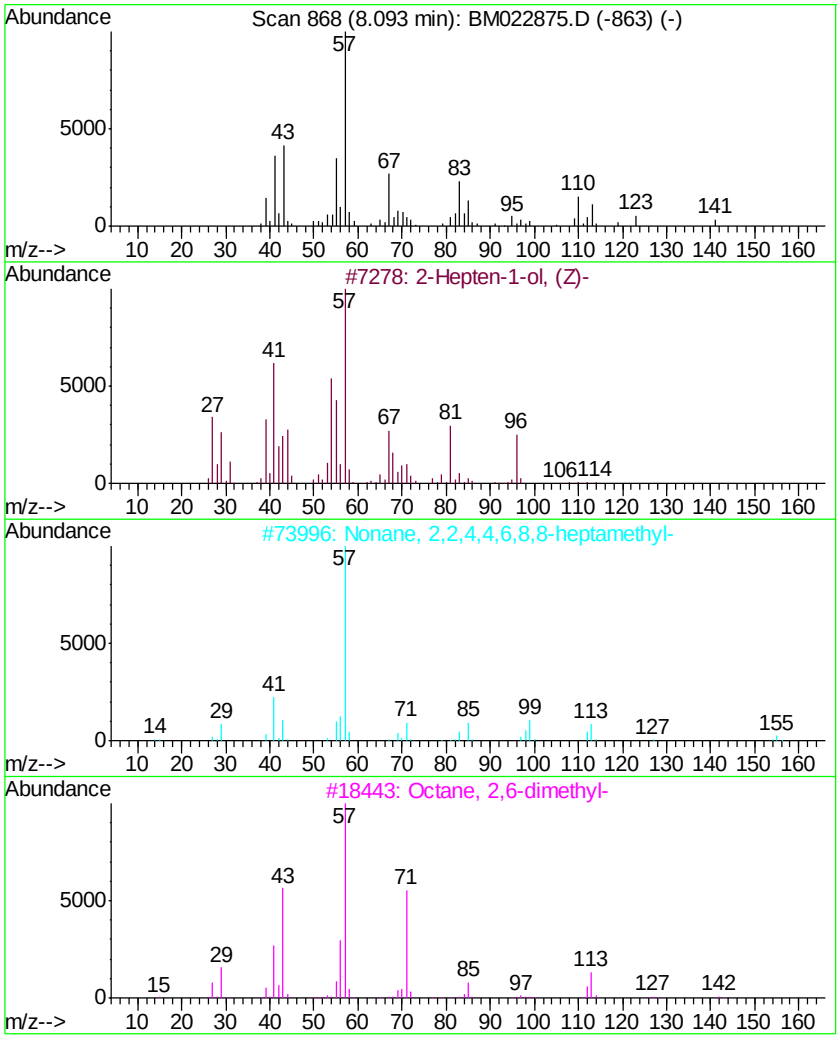
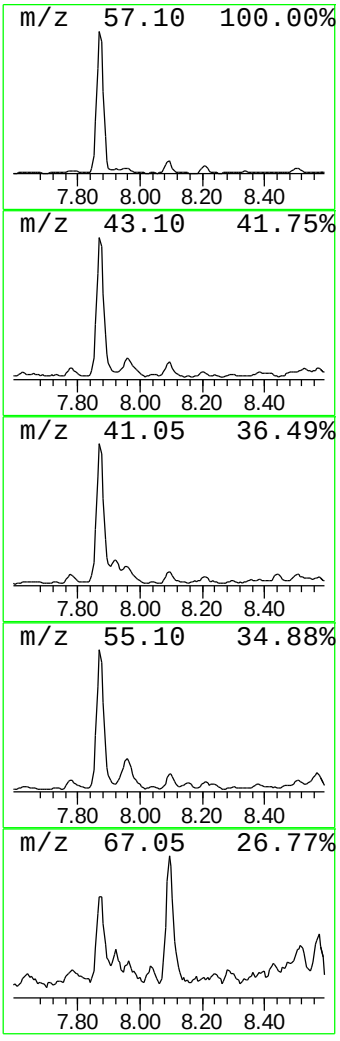
Peak Number 21 unknown-02 Concentration Rank 46

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.09	2.46 ng/ul	227153	1,4-Dichlorobenzene-d4	7.80

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Hepten-1-ol, (Z)-	114	C7H14O	055454-22-3	27
2	Nonane, 2,2,4,4,6,8,8-heptamethyl-	226	C16H34	004390-04-9	25
3	Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	25
4	Pentanoic acid, 2,4-dimethyl-, t...	186	C11H22O2	1000163-75-4	25
5	Nonane, 2,2,4,4,6,8,8-heptamethyl-	226	C16H34	004390-04-9	25





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Data File : BM022875.D
Acq On : 02 Oct 2019 04:21
Operator : JU
Sample : K4932-07
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BFDG9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM092719MA.M
Quant Title : SVOA CALIBRATION

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

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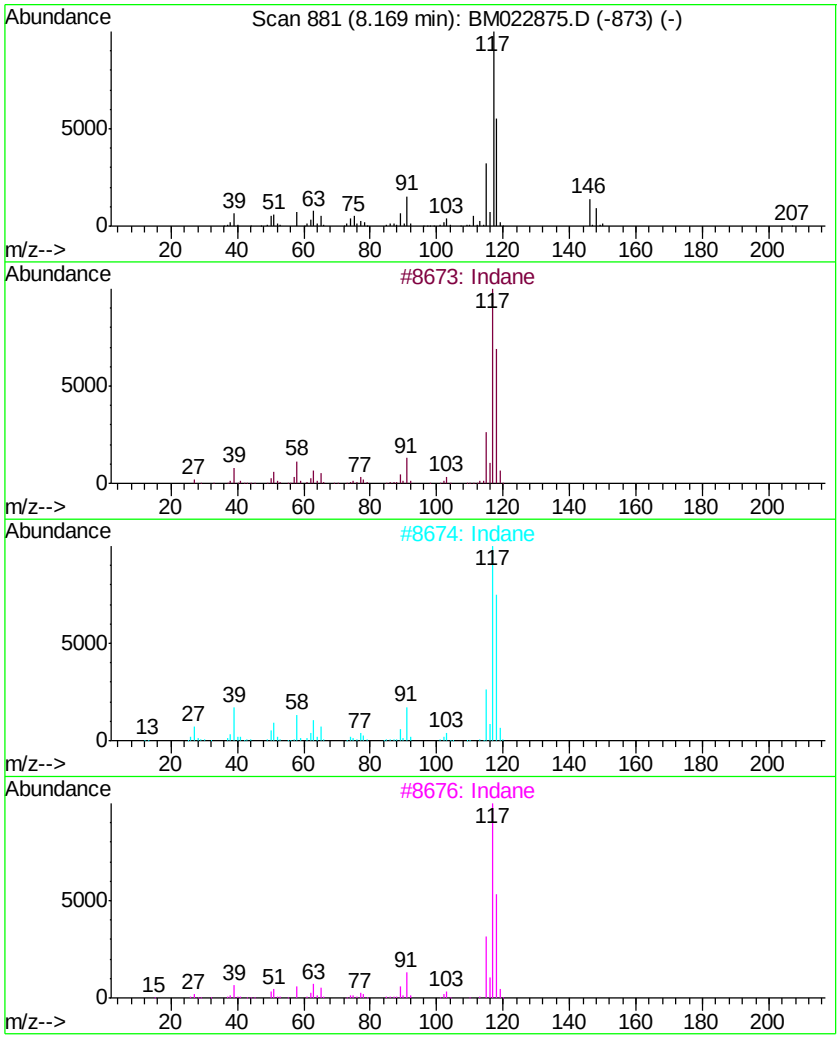
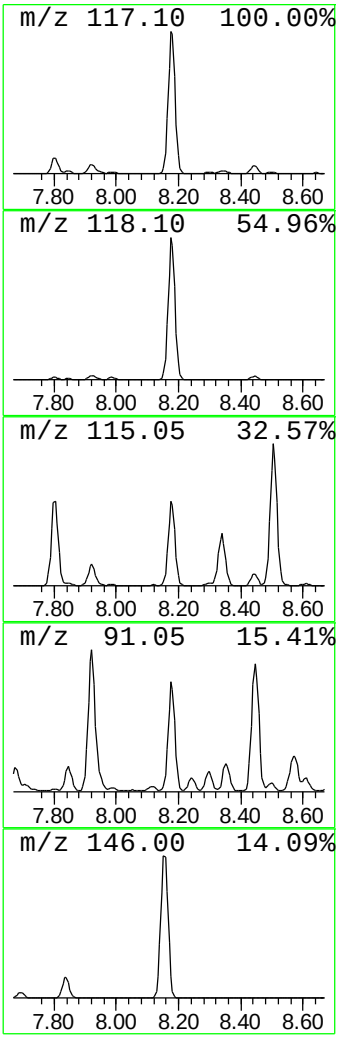
Peak Number 22 Indane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.17	20.90 ng/ul	1926130	1,4-Dichlorobenzene-d4	7.80

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane	118	C9H10	000496-11-7	81
2	Indane	118	C9H10	000496-11-7	81
3	Indane	118	C9H10	000496-11-7	72
4	Benzene, cyclopropyl-	118	C9H10	000873-49-4	72
5	cis-.beta.-Methylstyrene	118	C9H10	000766-90-5	68





Library Search Compound Report

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Acq On : 02 Oct 2019 04:21
Operator : JU
Sample : K4932-07
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BFDG9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM092719MA.M
Quant Title : SVOA CALIBRATION

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

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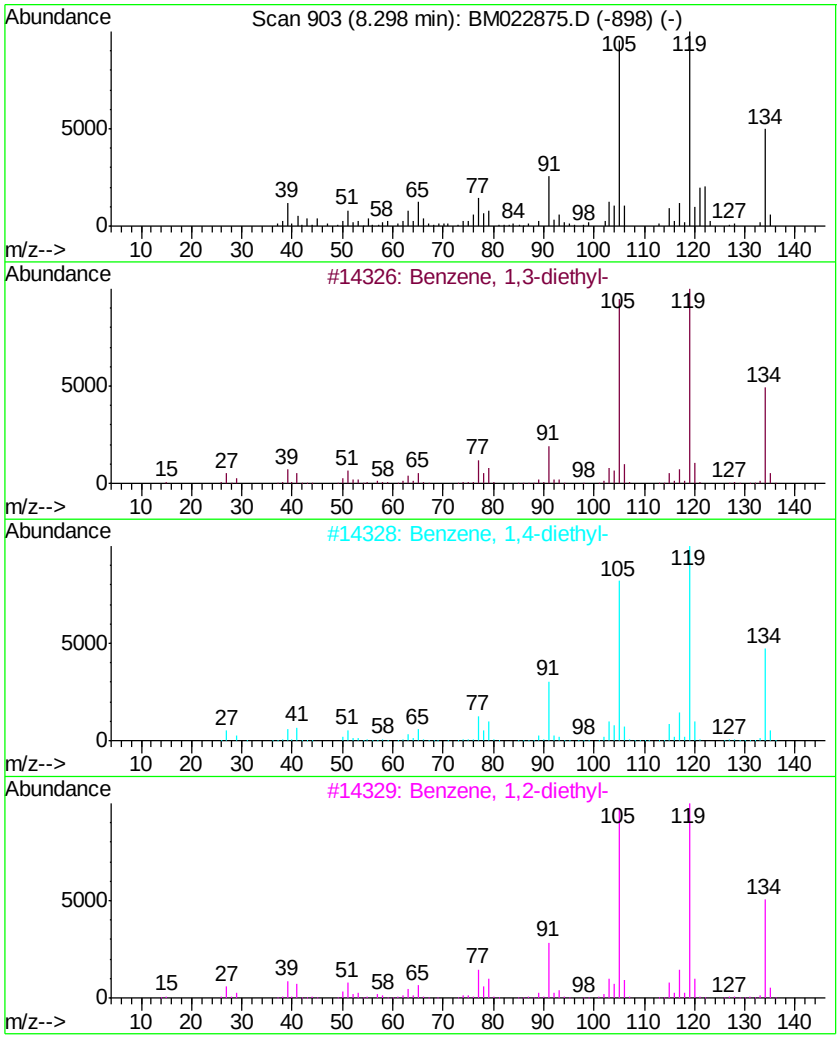
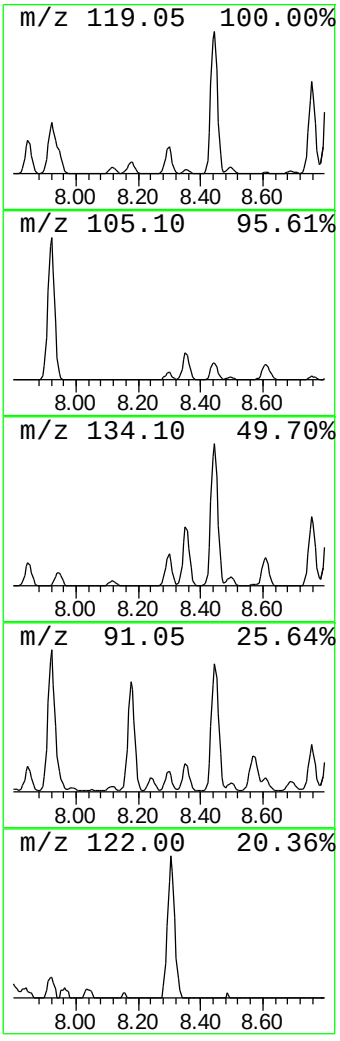
Peak Number 23 Benzene, 1,3-diethyl- Concentration Rank 35

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.30	3.18 ng/ul	293149	1,4-Dichlorobenzene-d4	7.80

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	96
2	Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	96
3	Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	96
4	Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	96
5	Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	93





Library Search Compound Report

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Data File : BM022875.D
Acq On : 02 Oct 2019 04:21
Operator : JU
Sample : K4932-07
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BFDG9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM092719MA.M
Quant Title : SVOA CALIBRATION

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

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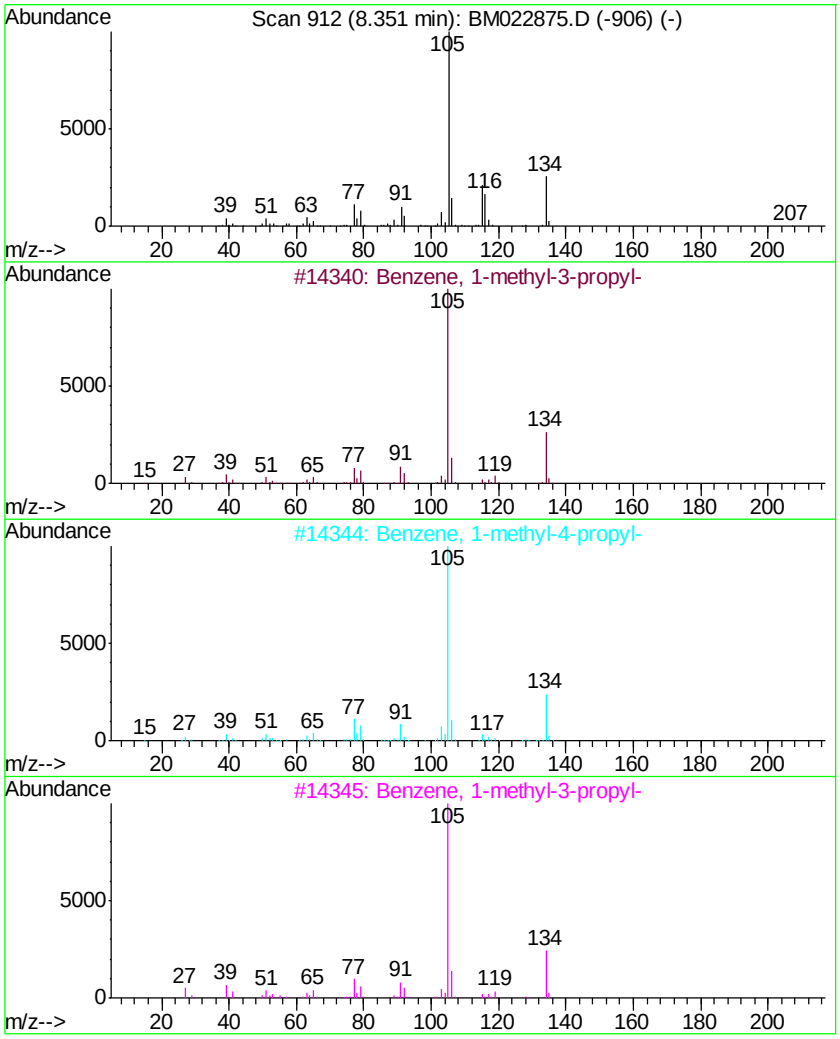
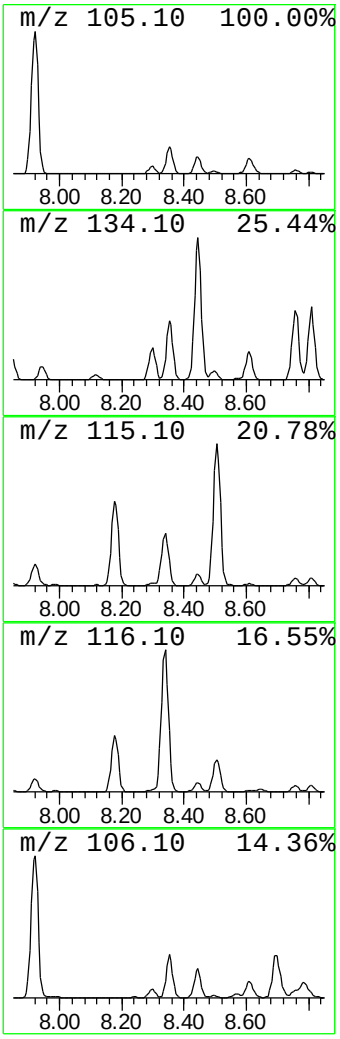
Peak Number 24 Benzene, 1-methyl-3-propyl- Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.35	7.75 ng/ul	714569	1,4-Dichlorobenzene-d4	7.80

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-methyl-3-propyl-	134	C10H14	001074-43-7	70
2	Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	70
3	Benzene, 1-methyl-3-propyl-	134	C10H14	001074-43-7	70
4	Benzene, 1-methyl-2-propyl-	134	C10H14	001074-17-5	62
5	2-Tolyloxirane	134	C9H10O	002783-26-8	58





Library Search Compound Report

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Acq On : 02 Oct 2019 04:21
Operator : JU
Sample : K4932-07
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BFDG9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM092719MA.M
Quant Title : SVOA CALIBRATION

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

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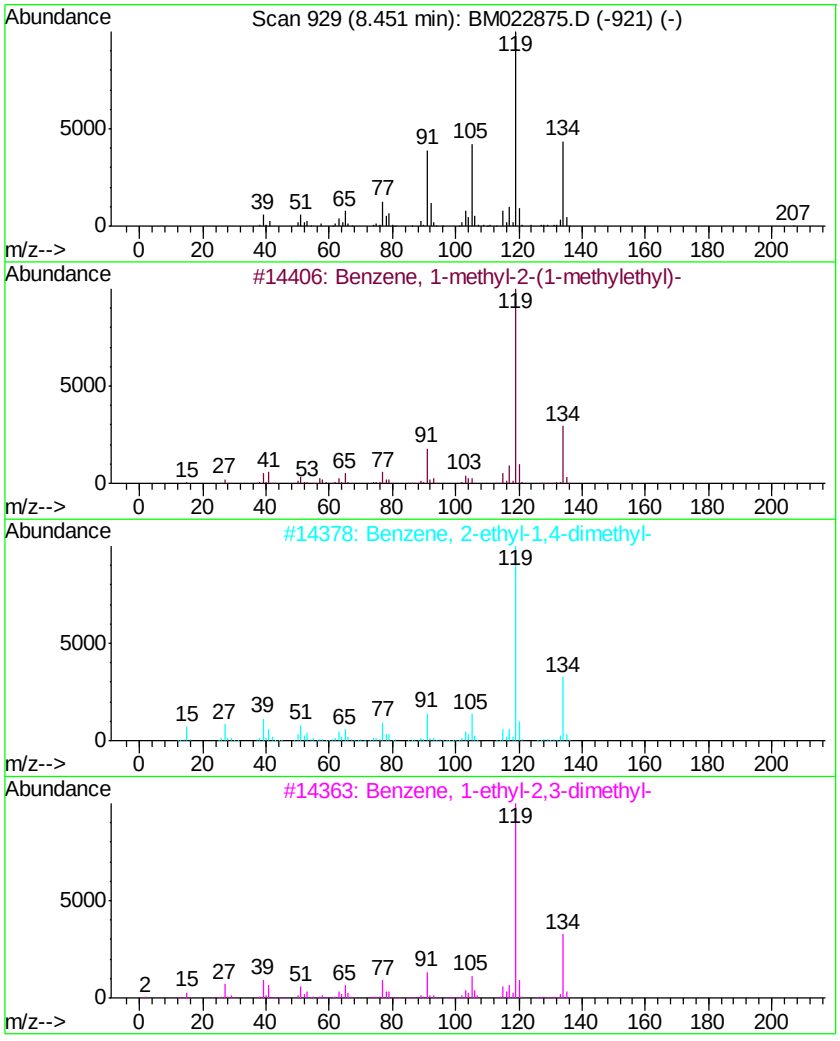
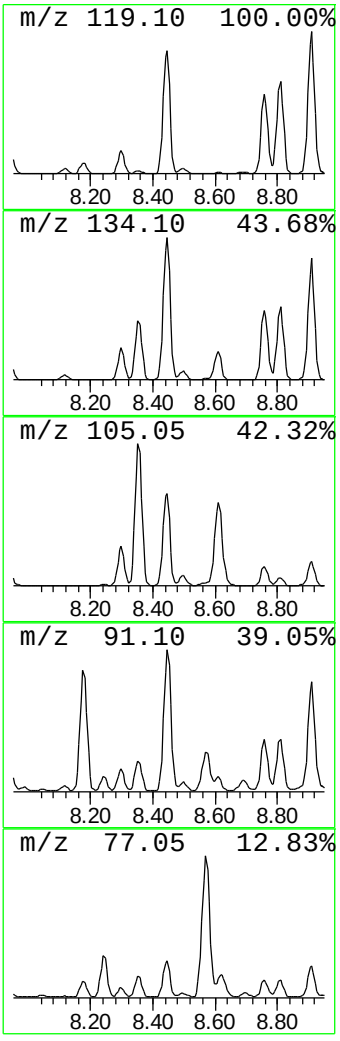
Peak Number 25 Benzene, 1-methyl-2-(1-meth... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.45	11.39 ng/ul	1049790	1,4-Dichlorobenzene-d4	7.80

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-methyl-2-(1-methyleth...	134	C10H14	000527-84-4	93
2	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	91
3	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	91
4	Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	91
5	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	91





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Acq On : 02 Oct 2019 04:21
Operator : JU
Sample : K4932-07
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BFDG9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM092719MA.M
Quant Title : SVOA CALIBRATION

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

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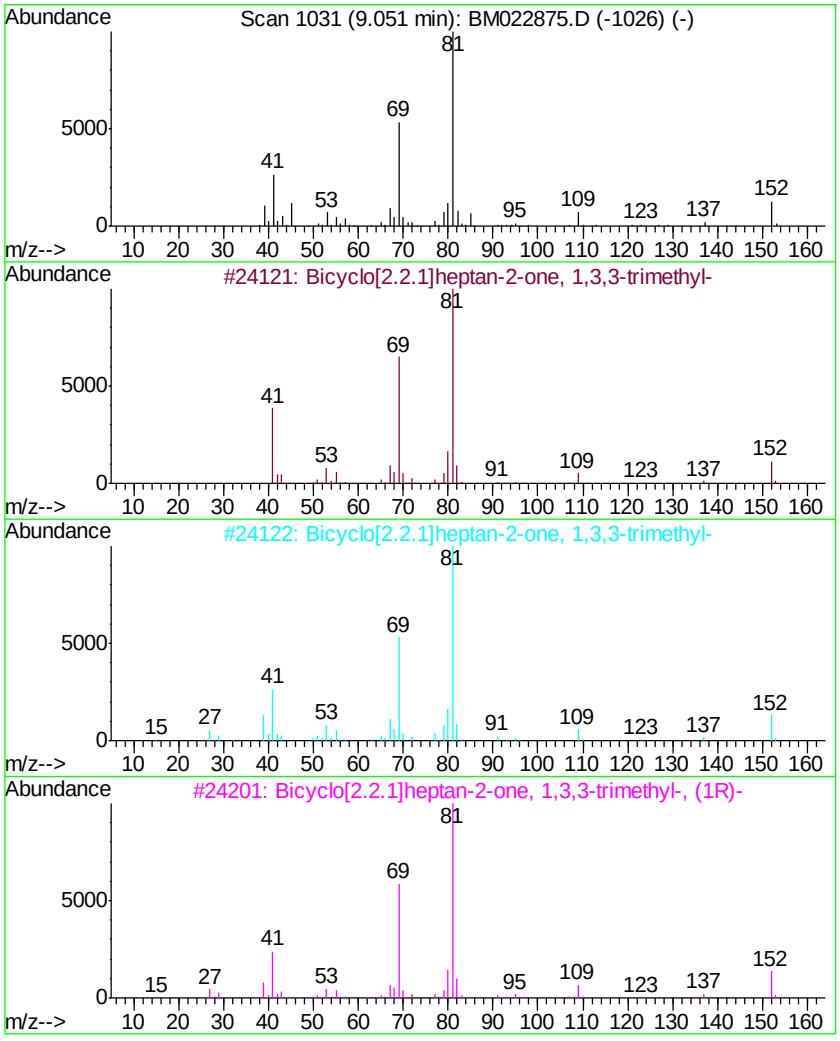
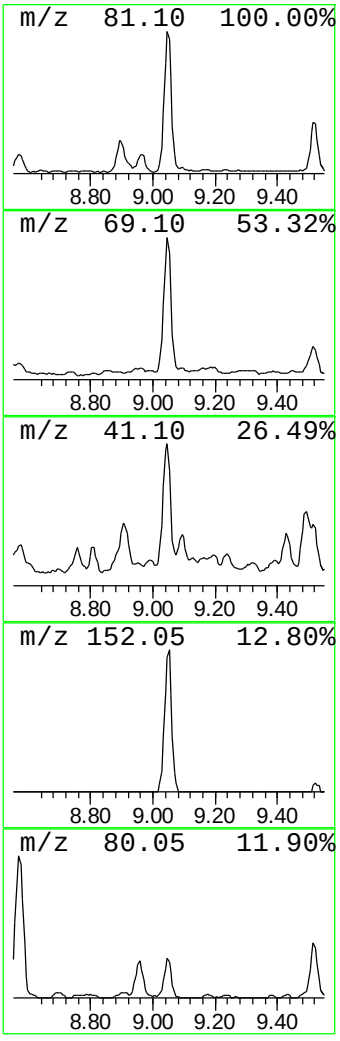
Peak Number 28 Bicyclo[2.2.1]heptan-2-one,... Concentration Rank 31

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.05	3.56 ng/ul	327651	1,4-Dichlorobenzene-d4	7.80

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Bicyclo[2.2.1]heptan-2-one, 1,3,...	152	C10H16O	001195-79-5	93
2	Bicyclo[2.2.1]heptan-2-one, 1,3,...	152	C10H16O	001195-79-5	93
3	Bicyclo[2.2.1]heptan-2-one, 1,3,...	152	C10H16O	007787-20-4	93
4	Bicyclo[2.2.1]heptan-2-one, 1,3,...	152	C10H16O	004695-62-9	76
5	L-Fenchone	152	C10H16O	000126-21-6	76





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM100119\
Data File : BM022875.D
Acq On : 02 Oct 2019 04:21
Operator : JU
Sample : K4932-07
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BFDG9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM092719MA.M
Quant Title : SVOA CALIBRATION

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

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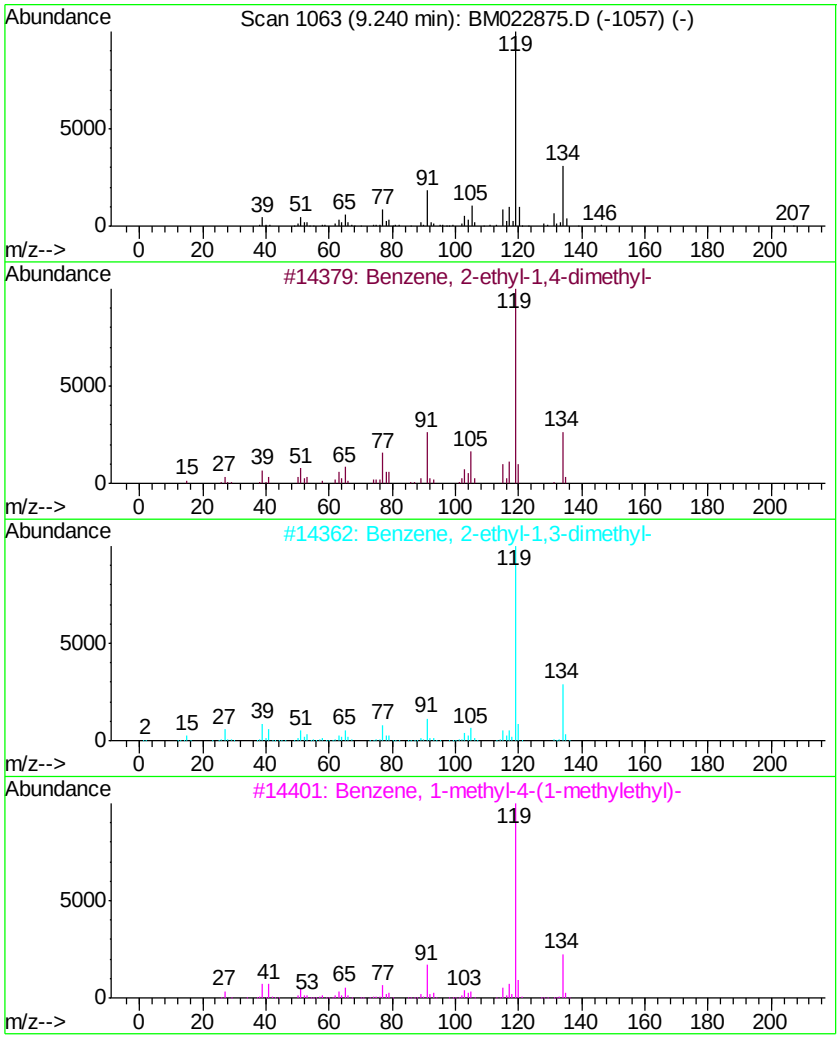
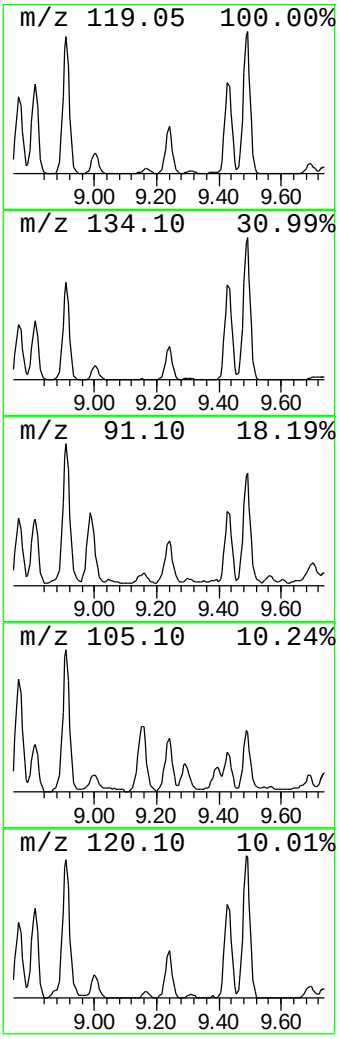
Peak Number 29 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 44

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.24	2.54 ng/ul	369903	Naphthalene-d8	10.59

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	96
2	Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	95
3	Benzene, 1-methyl-4-(1-methyleth...	134	C10H14	000099-87-6	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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Acq On : 02 Oct 2019 04:21
Operator : JU
Sample : K4932-07
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BFDG9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM092719MA.M
Quant Title : SVOA CALIBRATION

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

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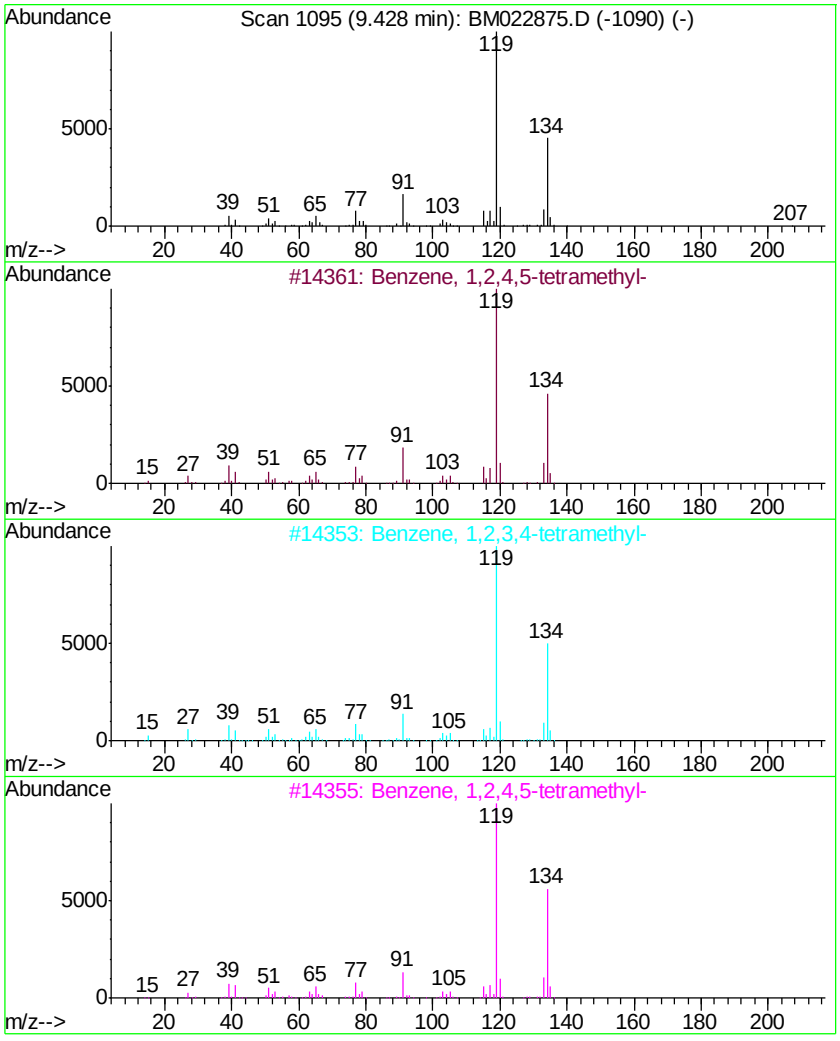
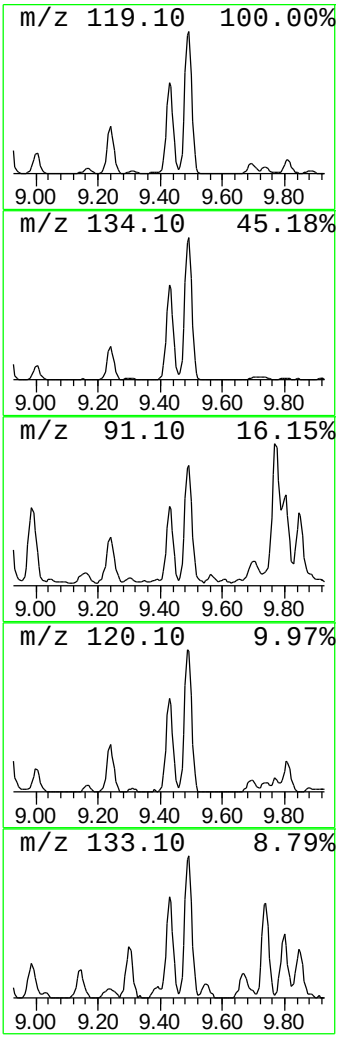
Peak Number 30 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 28

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.43	4.13 ng/ul	602559	Naphthalene-d8	10.59

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	95
2	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	95
3	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	95
4	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	95
5	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	94





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Acq On : 02 Oct 2019 04:21
Operator : JU
Sample : K4932-07
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BFDG9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM092719MA.M
Quant Title : SVOA CALIBRATION

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

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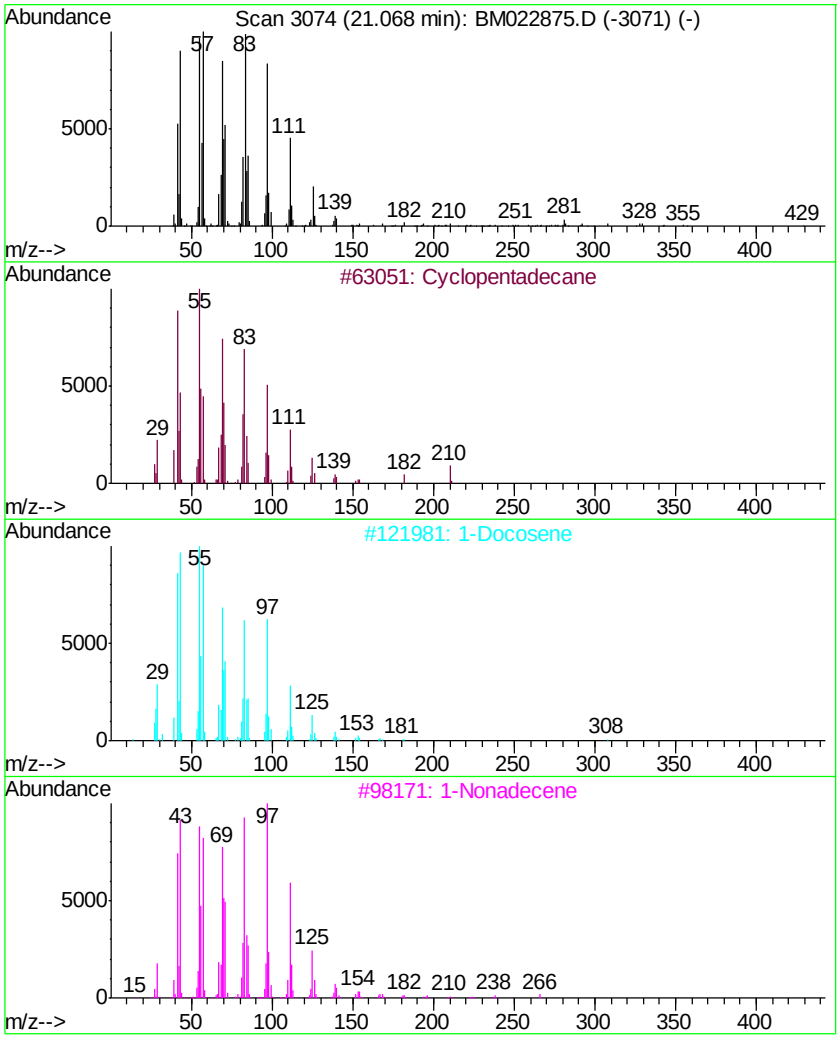
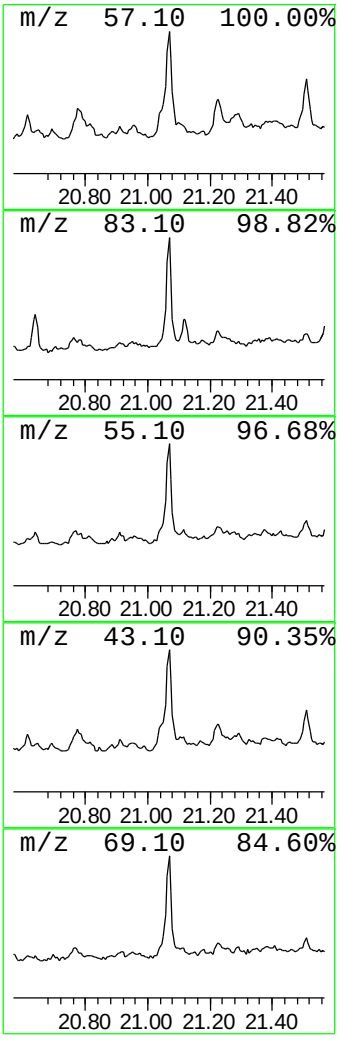
Peak Number 51 (DEL) Alkane: Cyclic21.07 Concentration Rank 32

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.07	3.51 ng/ul	514407	Chrysene-d12	21.36

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentadecane	210	C15H30	000295-48-7	93
2	1-Docosene	308	C22H44	001599-67-3	92
3	1-Nonadecene	266	C19H38	018435-45-5	91
4	Bromoacetic acid, octadecyl ester	390	C20H39BrO2	018992-03-5	91
5	3-Chloropropionic acid, heptadec...	346	C20H39ClO2	1000283-05-1	87





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ALS Vial : 21 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BFDG9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM092719MA.M
Quant Title : SVOA CALIBRATION

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

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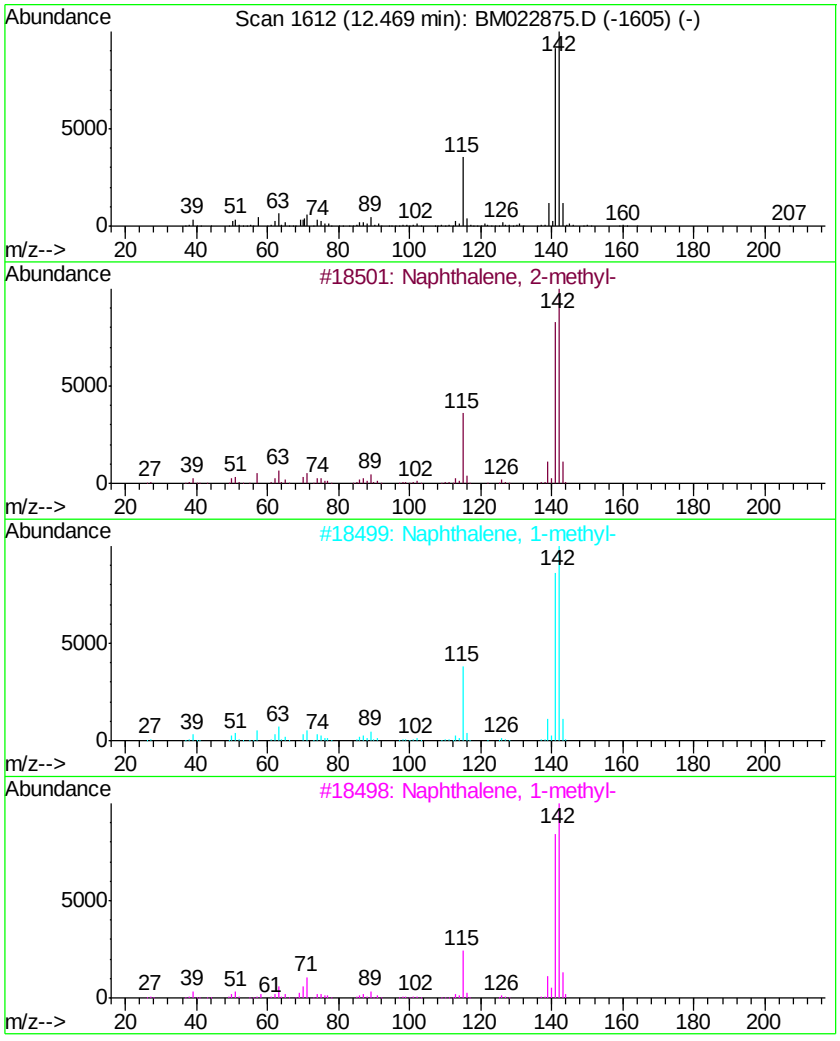
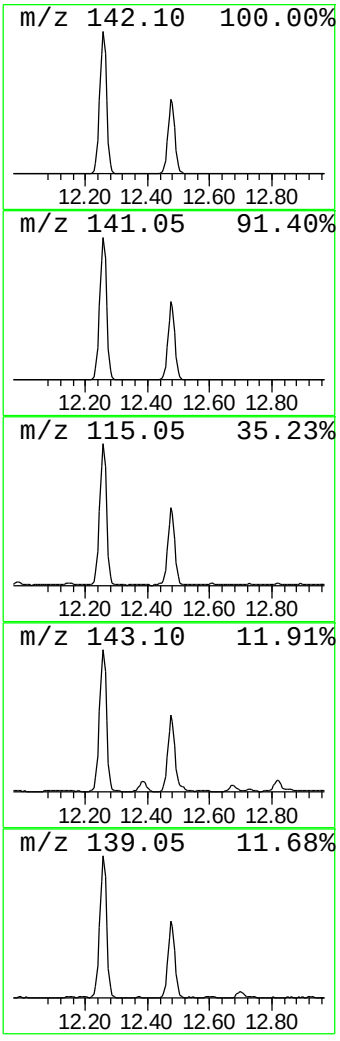
Peak Number 52 Naphthalene, 1-methyl- Concentration Rank 52

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.47	15.09 ng/ul	2200330	Naphthalene-d8	10.59

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
2	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
3	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	94
4	Benzocycloheptatriene	142	C11H10	000264-09-5	91
5	1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	91





Tentatively Identified Compound (LSC) summary

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Data File : BM022875.D
Acq On : 02 Oct 2019 04:21
Operator : JU
Sample : K4932-07
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BFDG9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM092719MA.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
Methyl Isobutyl K...	3.61	3.4	ng/ul	316358	1	7.80	1843180	20.0
Propanoic acid, 1...	3.73	3.1	ng/ul	285077	1	7.80	1843180	20.0
2-Hexanol	3.83	12.0	ng/ul	1102730	1	7.80	1843180	20.0
Toluene	4.00	79.9	ng/ul	7366880	1	7.80	1843180	20.0
Propanoic acid, 2...	4.27	6.5	ng/ul	602619	1	7.80	1843180	20.0
Propanoic acid, p...	4.45	10.0	ng/ul	925091	1	7.80	1843180	20.0
Butanoic acid, 1-...	4.90	12.2	ng/ul	1123020	1	7.80	1843180	20.0
Ethylbenzene	5.32	23.7	ng/ul	2188040	1	7.80	1843180	20.0
o-Xylene	5.46	90.4	ng/ul	8334880	1	7.80	1843180	20.0
p-Xylene	5.82	54.9	ng/ul	5062840	1	7.80	1843180	20.0
Benzene, (1-methy...	6.30	2.7	ng/ul	247418	1	7.80	1843180	20.0
Benzene, propyl-	6.79	7.4	ng/ul	684555	1	7.80	1843180	20.0
Benzene, 1-ethyl-...	6.90	26.8	ng/ul	2471950	1	7.80	1843180	20.0
unknown-01	7.26	2.8	ng/ul	255349	1	7.80	1843180	20.0
Benzene, 1,2,3-tr...	7.46	74.0	ng/ul	6822720	1	7.80	1843180	20.0
1-Hexanol, 2-ethyl-	7.87	14.7	ng/ul	1351970	1	7.80	1843180	20.0
unknown-02	8.09	2.5	ng/ul	227153	1	7.80	1843180	20.0
Indane	8.17	20.9	ng/ul	1926130	1	7.80	1843180	20.0
Benzene, 1,3-diet...	8.30	3.2	ng/ul	293149	1	7.80	1843180	20.0
Benzene, 1-methyl...	8.35	7.8	ng/ul	714569	1	7.80	1843180	20.0
Benzene, 1-methyl...	8.45	11.4	ng/ul	1049790	1	7.80	1843180	20.0
Bicyclo[2.2.1]hep...	9.05	3.6	ng/ul	327651	1	7.80	1843180	20.0
Benzene, 2-ethyl-...	9.24	2.5	ng/ul	369903	2	10.59	2916970	20.0
Benzene, 1,2,4,5-...	9.43	4.1	ng/ul	602559	2	10.59	2916970	20.0
(DEL) Alkane: Cyc...	21.07	3.5	ng/ul	514407	5	21.36	2931360	20.0
Naphthalene, 1-me...	12.47	15.1	ng/ul	2200330	2	10.59	2916970	20.0