LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM060420\
Data File : BM026264.D
Acq On : 04 Jun 2020 19:12
Operator : JU/CG
Sample : L2844-15
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C0AF2

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-BM051320MA.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	3.052	25	28	35	rBV	48214	62666	1.74%	0.169%
2	3.293	60	69	73	rBV2	35671	72283	2.01%	0.194%
3	3.434	90	93	101	rVV	15200	23772	0.66%	0.064%
4	3.516	103	107	114	rVB2	17140	26439	0.73%	0.071%
5	3.669	126	133	139	rVB3	17614	36679	1.02%	0.099%
6	3.752	140	147	158	rVB4	22611	56803	1.58%	0.153%
7	3.922	170	176	179	rBV2	38123	59371	1.65%	0.160%
8	3.958	179	182	188	rVB2	21134	30113	0.84%	0.081%
9	4.052	192	198	202	rVV6	17498	32670	0.91%	0.088%
10	4.193	218	222	230	rVB2	41621	59583	1.65%	0.160%
11	4.269	230	235	242	rBV2	21852	35506	0.99%	0.096%
12	4.628	291	296	299	rVV2	17134	27704	0.77%	0.075%
13	4.669	299	303	308	rVV	41949	58264	1.62%	0.157%
14	4.722	308	312	318	rVV	37409	55288	1.53%	0.149%
15	4.958	343	352	353	rBV4	14907	27029	0.75%	0.073%
16	5.152	380	385	392	rBV	31417	66352	1.84%	0.178%
17	5.210	392	395	402	rVB2	16798	26630	0.74%	0.072%
18	5.381	420	424	430	rBV5	15586	26644	0.74%	0.072%
19	5.457	432	437	442	rVV3	26400	46249	1.28%	0.124%
20	5.522	442	448	455	rVV3	36634	68815	1.91%	0.185%
21	5.775	484	491	503	rBV5	29652	83016	2.30%	0.223%
22	5.993	518	528	534	rBV3	44978	87358	2.43%	0.235%
23	6.122	543	550	555	rBV3	32804	62781	1.74%	0.169%
24	6.463	601	608	615	rBV5	25750	51559	1.43%	0.139%
25	6.575	623	627	631	rVV	49925	71185	1.98%	0.191%
26	6.634	631	637	638	rVV2	29270	59045	1.64%	0.159%
27	6.663	638	642	646	rVV	177592	285121	7.91%	0.767%
28	6.704	646	649	654	rVB	113465	158099	4.39%	0.425%
29	6.816	661	668	674	rBV	540844	833522	23.14%	2.242%
30	6.863	674	676	681	rVB2	29762	38603	1.07%	0.104%
31	7.004	694	700	708	rVV	606154	931052	25.85%	2.505%
32	7.122	714	720	725	rBV	153126	223742	6.21%	0.602%
33	7.463	771	778	783	rBV	530512	785069	21.79%	2.112%
34	7.510	783	786	790	rVV2	26181	32997	0.92%	0.089%

LSC Area Percent Report

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Operator : JU/CG
Sample : L2844-15
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C0AF2

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-BM051320MA.M
Title : SVOA CALIBRATION

35	7.581	790	798	805	rVB2	52355	101436	2.82%	0.273%
36	7.957	857	862	867	rBV3	14733	25213	0.70%	0.068%
37	8.010	867	871	877	rVV	45386	67476	1.87%	0.182%
38	8.104	881	887	892	rVV	84044	130052	3.61%	0.350%
39	8.181	892	900	910	rVV	458249	731165	20.30%	1.967%
40	8.269	910	915	922	rVV2	39097	82629	2.29%	0.222%
41	8.410	934	939	944	rBV2	27141	41491	1.15%	0.112%
42	8.463	944	948	954	rVB	34612	48180	1.34%	0.130%
43	8.563	956	965	967	rBV	75772	116772	3.24%	0.314%
44	8.604	967	972	984	rVV	669483	1109020	30.79%	2.983%
45	8.704	984	989	995	rVB3	15010	28001	0.78%	0.075%
46	8.810	995	1007	1013	rVB4	38877	84003	2.33%	0.226%
47	8.893	1015	1021	1026	rBV2	25379	46728	1.30%	0.126%
48	8.957	1026	1032	1037	rVB3	27834	49296	1.37%	0.133%
49	9.081	1048	1053	1058	rVB2	58864	89163	2.48%	0.240%
50	9.140	1058	1063	1069	rVB	80309	129041	3.58%	0.347%
51	9.322	1087	1094	1101	rBV	572635	899859	24.98%	2.421%
52	9.451	1110	1116	1125	rVB4	35638	82326	2.29%	0.221%
53	9.645	1143	1149	1154	rVB2	79175	133542	3.71%	0.359%
54	9.710	1156	1160	1166	rVB3	28434	48011	1.33%	0.129%
55	9.798	1170	1175	1178	rBV2	24356	42509	1.18%	0.114%
56	9.857	1178	1185	1195	rVB	740516	1216802	33.78%	3.273%
57	9.945	1195	1200	1205	rVB	28680	51550	1.43%	0.139%
58	10.128	1223	1231	1232	rBV6	13314	25413	0.71%	0.068%
59	10.151	1232	1235	1240	rVB5	16134	24490	0.68%	0.066%
60	10.222	1240	1247	1253	rBV	682938	1126957	31.28%	3.032%
61	10.287	1253	1258	1261	rBV3	25427	44724	1.24%	0.120%
62	10.375	1270	1273	1278	rVB2	33326	48519	1.35%	0.131%
63	10.457	1283	1287	1295	rVB2	19765	34822	0.97%	0.094%
64	10.604	1306	1312	1317	rBV3	17211	26872	0.75%	0.072%
65	10.687	1317	1326	1332	rBV6	18041	44502	1.24%	0.120%
66	11.287	1422	1428	1433	rBV7	10808	24711	0.69%	0.066%
67	11.416	1442	1450	1455	rBV5	17492	33954	0.94%	0.091%
68	11.781	1508	1512	1516	rVV3	29532	52633	1.46%	0.142%
69	11.892	1529	1531	1537	rVB4	18766	31389	0.87%	0.084%
70	12.116	1565	1569	1574	rVV	34177	53276	1.48%	0.143%
71	13.045	1720	1727	1734	rVB7	16264	33531	0.93%	0.090%

LSC Area Percent Report

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Operator : JU/CG
Sample : L2844-15
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C0AF2

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-BM051320MA.M
Title : SVOA CALIBRATION

72	13.269	1760	1765	1766	rBV	18615	25167	0.70%	0.068%
73	13.428	1782	1792	1797	rVV2	44743	89039	2.47%	0.240%
74	13.480	1797	1801	1804	rVV2	24530	34985	0.97%	0.094%
75	13.533	1804	1810	1815	rVV	1386218	1849418	51.34%	4.975%
76	13.580	1815	1818	1823	rVV	178803	235884	6.55%	0.635%
77	13.663	1828	1832	1835	rVV2	26998	38230	1.06%	0.103%
78	13.698	1835	1838	1843	rVB3	18252	25434	0.71%	0.068%
79	13.798	1848	1855	1865	rVV	1486393	2144654	59.53%	5.769%
80	14.110	1900	1908	1916	rBV	917303	1337139	37.12%	3.597%
81	14.345	1939	1948	1957	rBV3	101022	223417	6.20%	0.601%
82	14.522	1968	1978	1985	rVV6	25411	75775	2.10%	0.204%
83	14.598	1987	1991	1996	rVV	25307	38629	1.07%	0.104%
84	14.739	2010	2015	2021	rVV3	27879	49961	1.39%	0.134%
85	14.916	2037	2045	2047	rBV4	13961	30309	0.84%	0.082%
86	15.110	2066	2078	2084	rBV	1975889	2737469	75.99%	7.364%
87	15.163	2084	2087	2093	rVV4	26316	42763	1.19%	0.115%
88	15.245	2093	2101	2113	rVV	670238	976446	27.11%	2.627%
89	15.351	2115	2119	2123	rVV2	26235	43840	1.22%	0.118%
90	15.898	2209	2212	2216	rVV3	18391	25515	0.71%	0.069%
91	16.669	2337	2343	2348	rVV6	10768	28457	0.79%	0.077%
92	16.863	2369	2376	2381	rBV	1084777	1526925	42.39%	4.108%
93	16.957	2386	2392	2406	rVV	2156861	3039151	84.37%	8.176%
94	18.898	2718	2722	2728	rVB	25969	35694	0.99%	0.096%
95	19.151	2760	2765	2769	rVB3	23367	33749	0.94%	0.091%
96	19.262	2778	2784	2791	rBV	2756476	3602360	100.00%	9.691%
97	20.809	3043	3047	3053	rBV	170938	248583	6.90%	0.669%
98	21.062	3085	3090	3098	rBV	1505121	1829104	50.78%	4.921%
99	23.045	3421	3427	3433	rVB	2022313	3323600	92.26%	8.941%
100	23.174	3443	3449	3459	rVB	1087789	1918923	53.27%	5.162%

Sum of corrected areas: 37172687

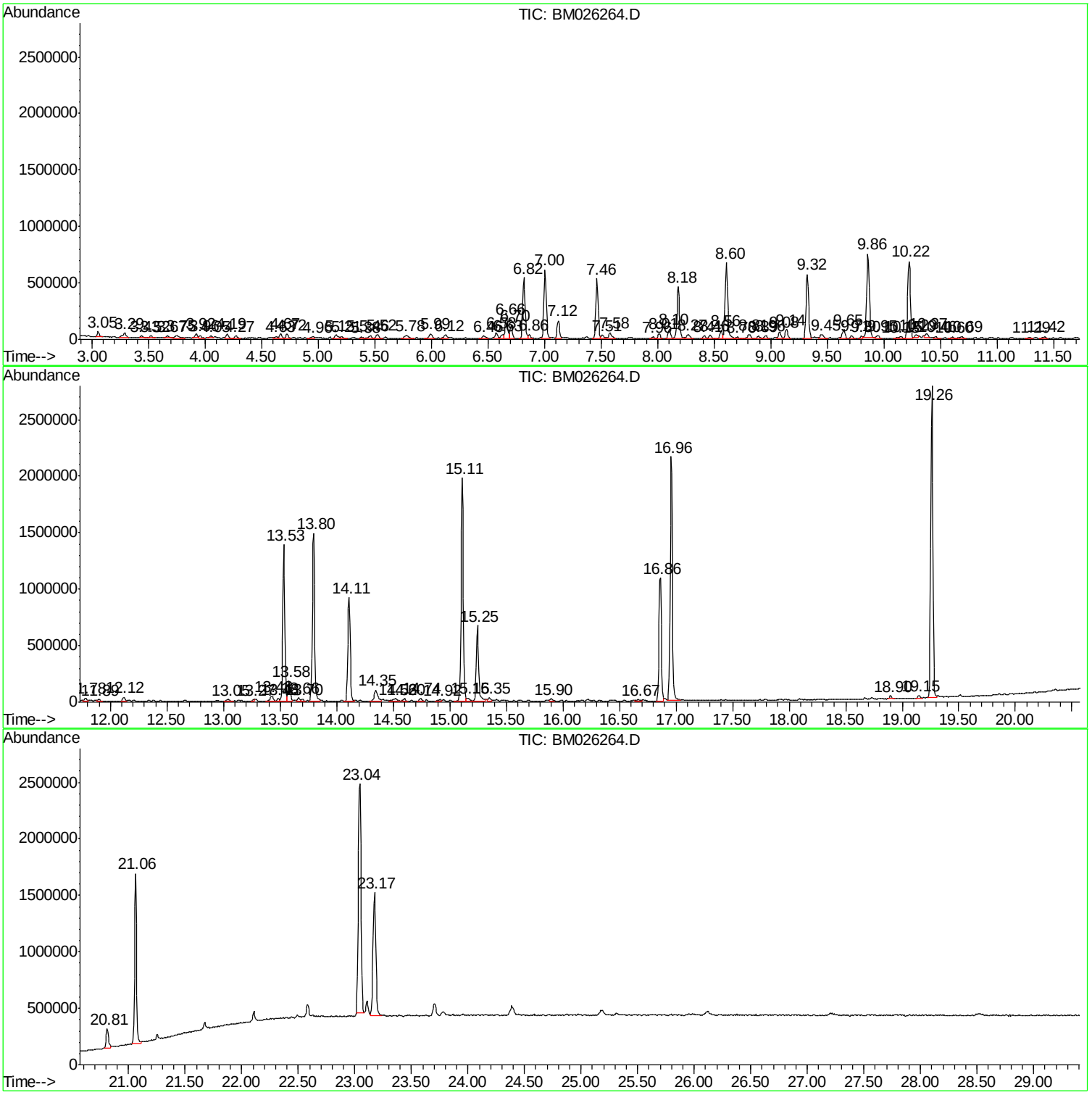
LSC Report - Integrated Chromatogram

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM060420\
Data File : BM026264.D
Acq On : 04 Jun 2020 19:12
Operator : JU/CG
Sample : L2844-15
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C0AF2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM060420\
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Instrument : BNA_M
ClientSampleId : C0AF2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION

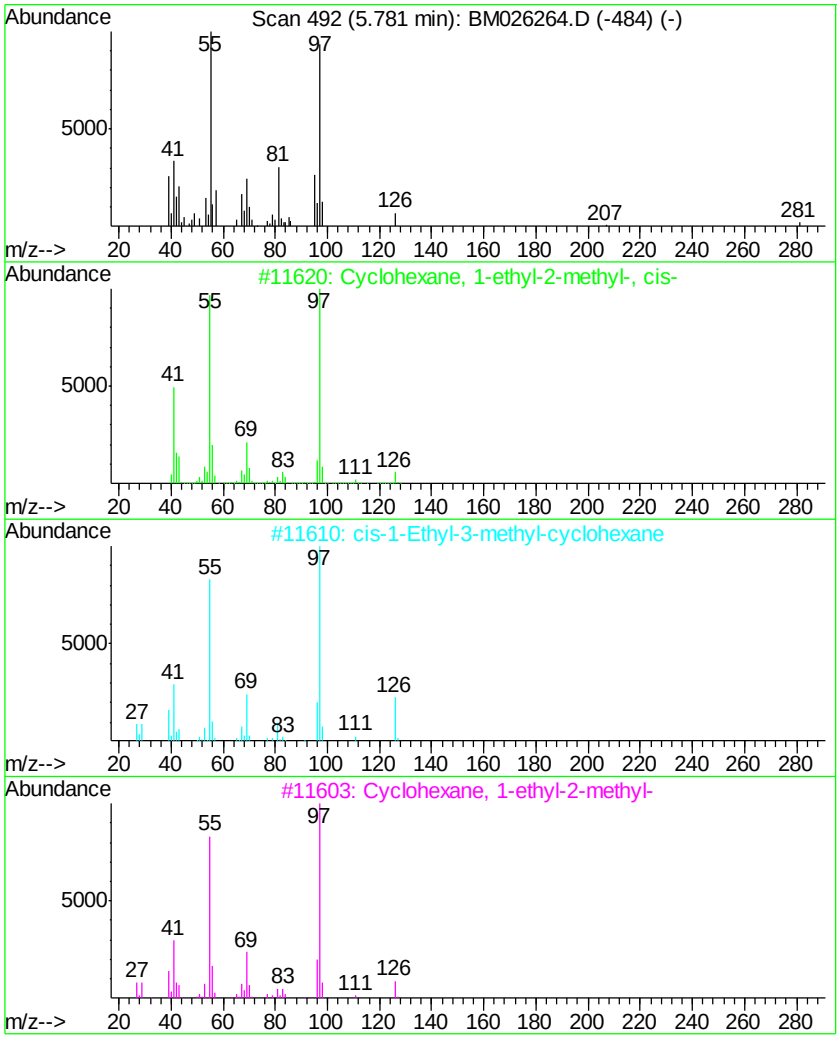
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

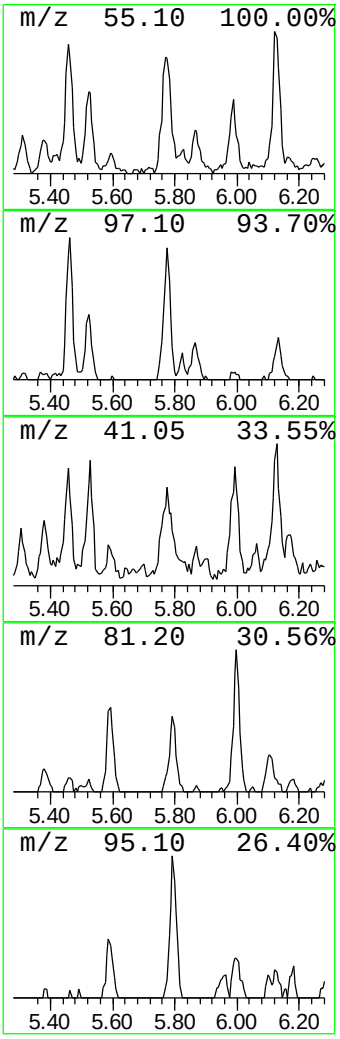
****

Peak Number 1 (DEL) Alkane: Cyclic5.78 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.78	2.11 ng/ul	83016	1,4-Dichlorobenzene-d4	7.46

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexane, 1-ethyl-2-methyl-, ...	126	C9H18	004923-77-7	87
2	cis-1-Ethyl-3-methyl-cyclohexane	126	C9H18	019489-10-2	83
3	Cyclohexane, 1-ethyl-2-methyl-	126	C9H18	003728-54-9	81
4	Cyclohexane, 1-ethyl-2-methyl-, ...	126	C9H18	004923-78-8	81
5	Cyclohexane, 1-ethyl-4-methyl-, ...	126	C9H18	004926-78-7	81





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM060420\
Data File : BM026264.D
Acq On : 04 Jun 2020 19:12
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Sample : L2844-15
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Instrument : BNA_M
ClientSampleId : C0AF2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION

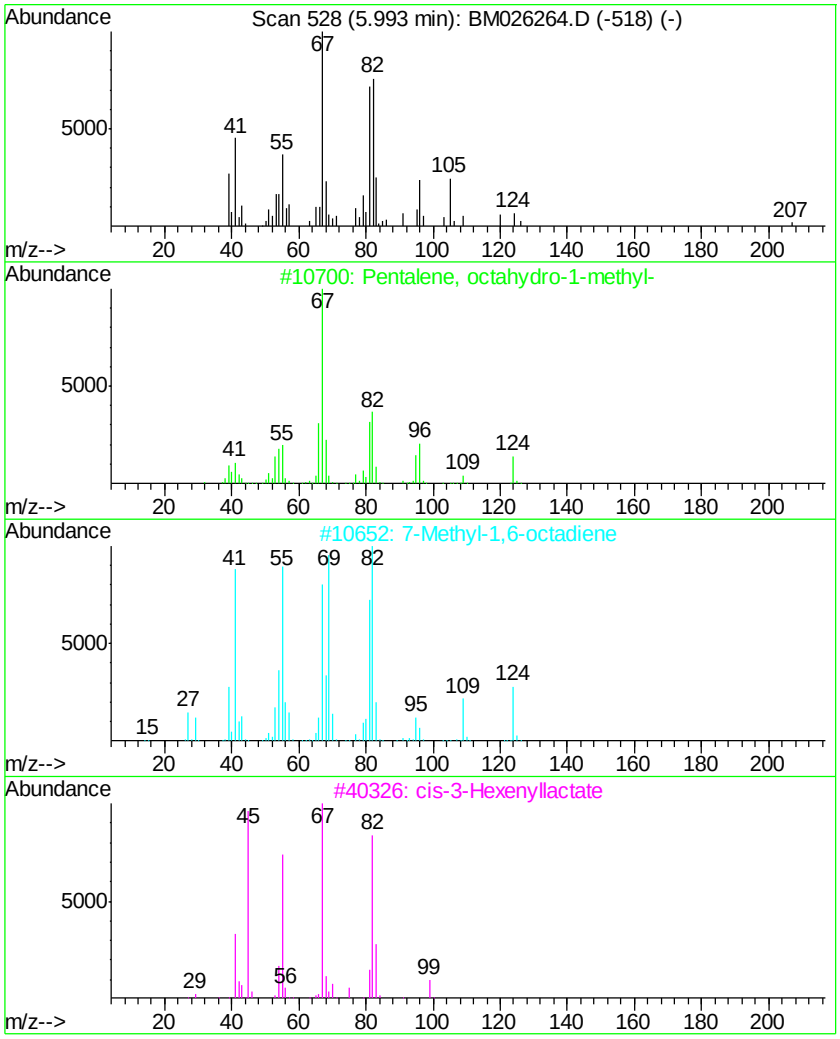
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

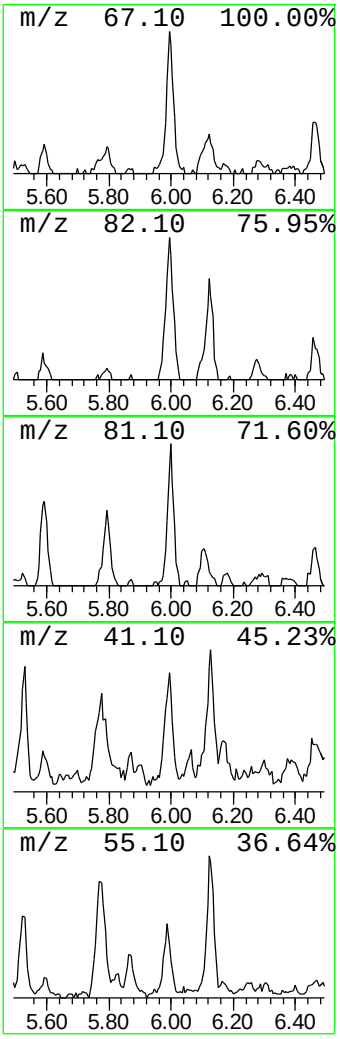
****

Peak Number 2 Pentalene, octahydro-1-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.99	2.23 ng/ul	87358	1,4-Dichlorobenzene-d4	7.46

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentalene, octahydro-1-methyl-	124	C9H16	032273-77-1	64
2	7-Methyl-1,6-octadiene	124	C9H16	042152-47-6	58
3	cis-3-Hexenyllactate	172	C9H16O3	061931-81-5	53
4	2,4-Hexadiene, (E,Z)-	82	C6H10	005194-50-3	52
5	1,3-Pentadiene, 2-methyl-, (E)-	82	C6H10	000926-54-5	52





Library Search Compound Report

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION

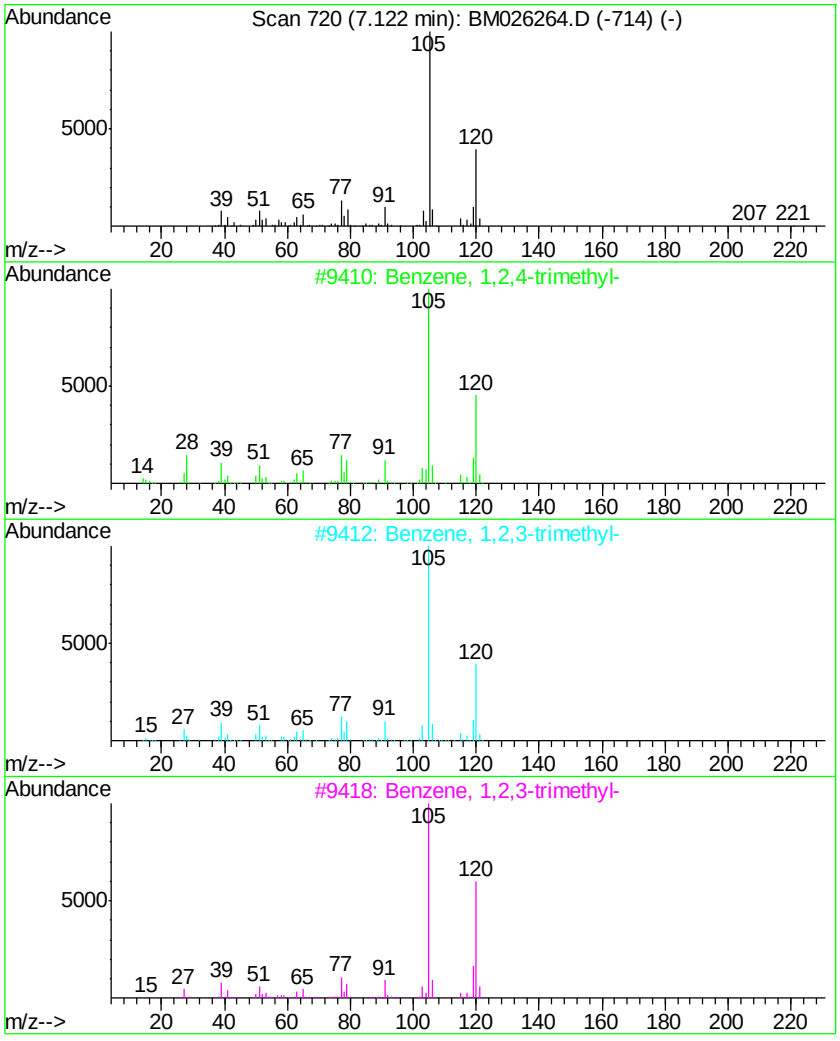
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

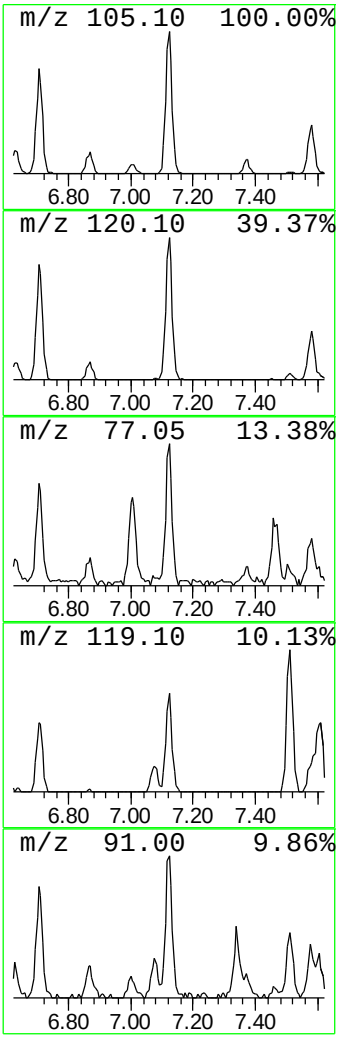
****

Peak Number 3 Benzene, 1,2,4-trimethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.12	5.70 ng/ul	223742	1,4-Dichlorobenzene-d4	7.46

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94
2	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	94
3	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	94
4	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	93
5	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	93





Library Search Compound Report

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Instrument : BNA_M
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Quant Title : SVOA CALIBRATION

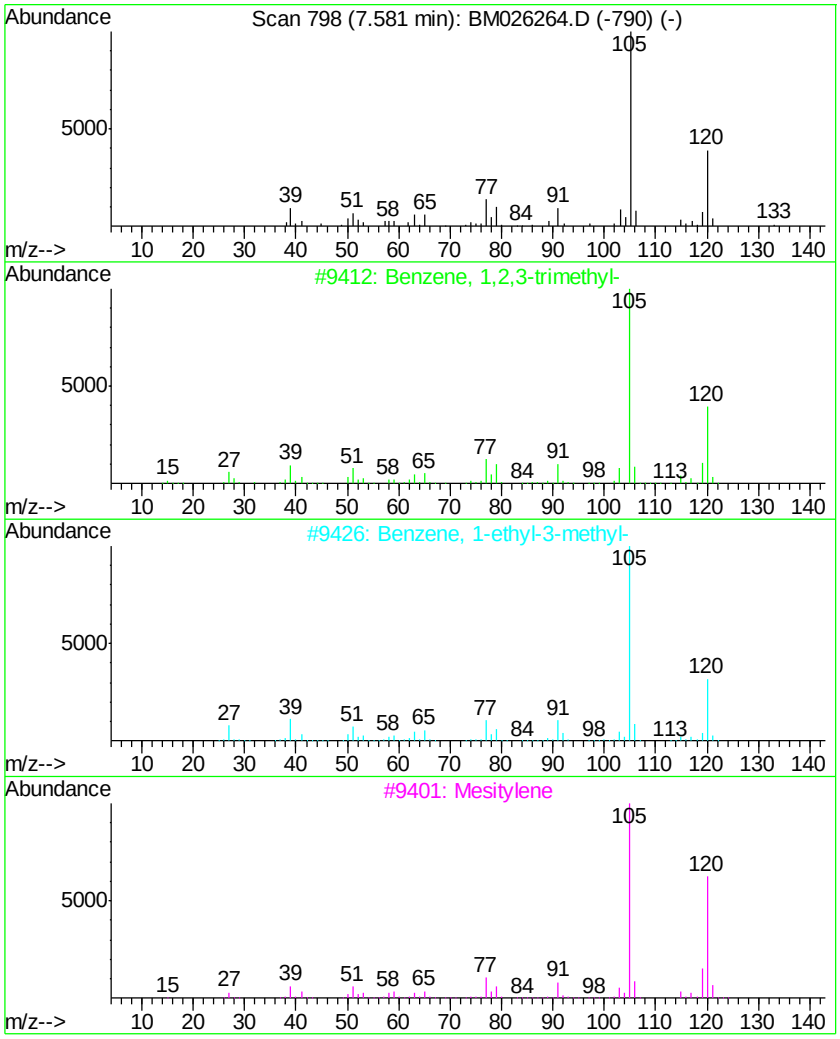
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

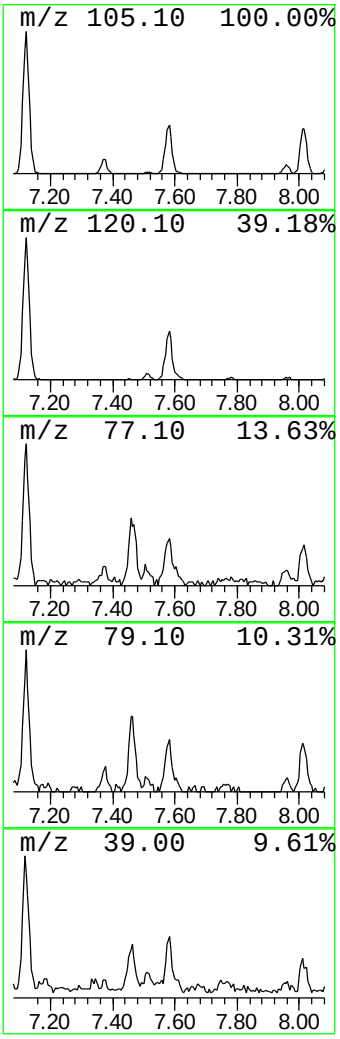
****

Peak Number 4 Benzene, 1,2,3-trimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.58	2.58 ng/ul	101436	1,4-Dichlorobenzene-d4	7.46

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
2	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
3	Mesitylene	120	C9H12	000108-67-8	91
4	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
5	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91





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Data File : BM026264.D
Acq On : 04 Jun 2020 19:12
Operator : JU/CG
Sample : L2844-15
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C0AF2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION

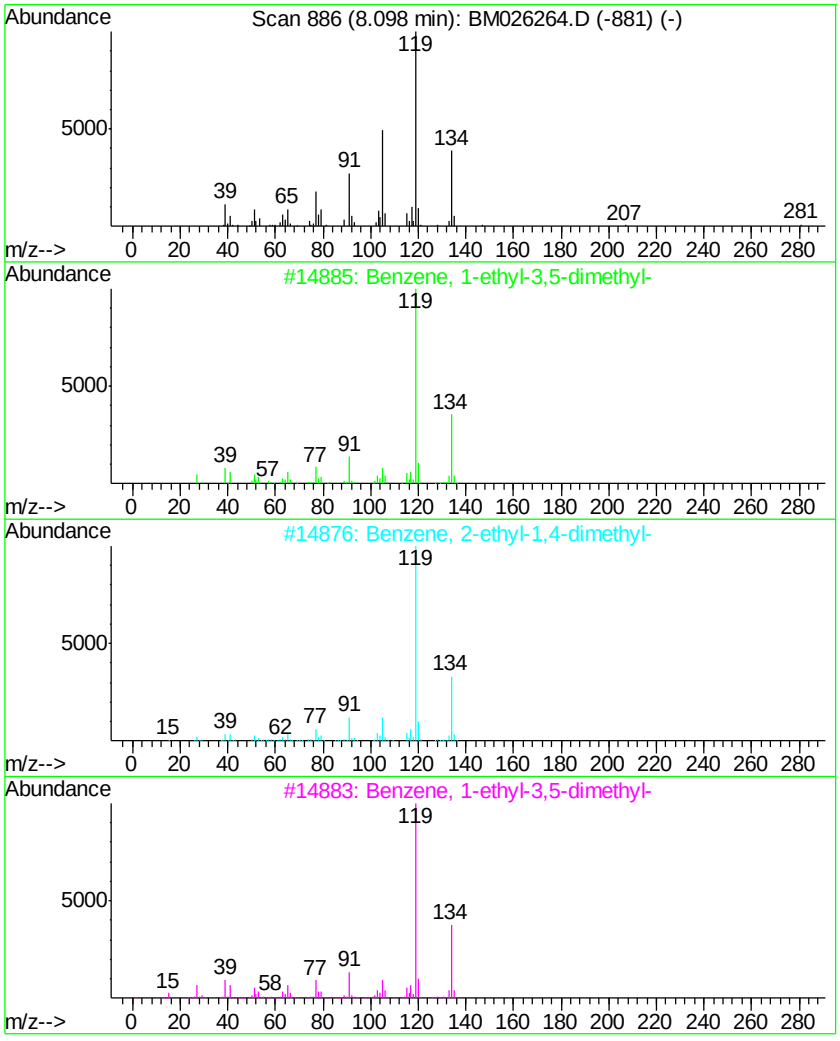
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

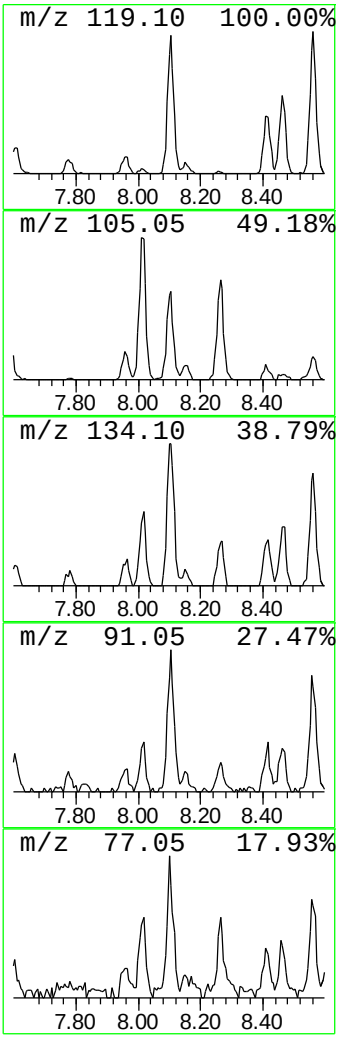
****

Peak Number 5 Benzene, 1-ethyl-3,5-dimethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.10	3.31 ng/ul	130052	1,4-Dichlorobenzene-d4	7.46

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	95
2	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	95
3	Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	95
4	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	94
5	p-Cymene	134	C10H14	000099-87-6	94





Library Search Compound Report

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Data File : BM026264.D
Acq On : 04 Jun 2020 19:12
Operator : JU/CG
Sample : L2844-15
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C0AF2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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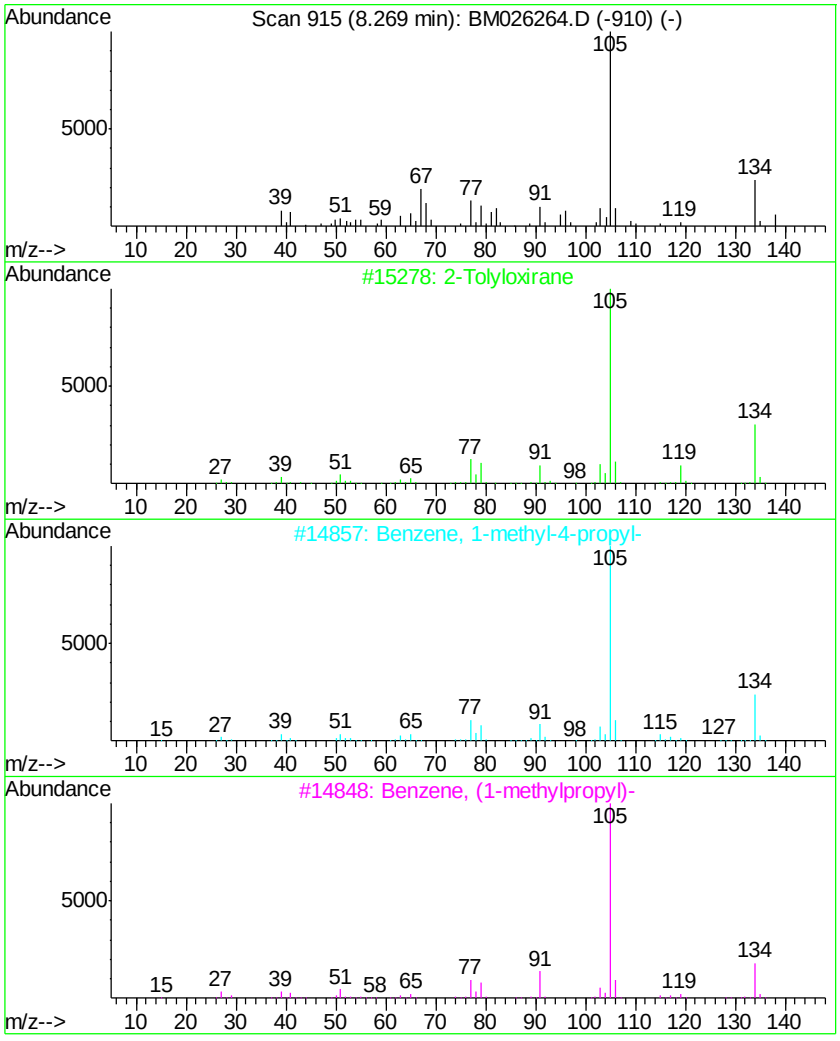
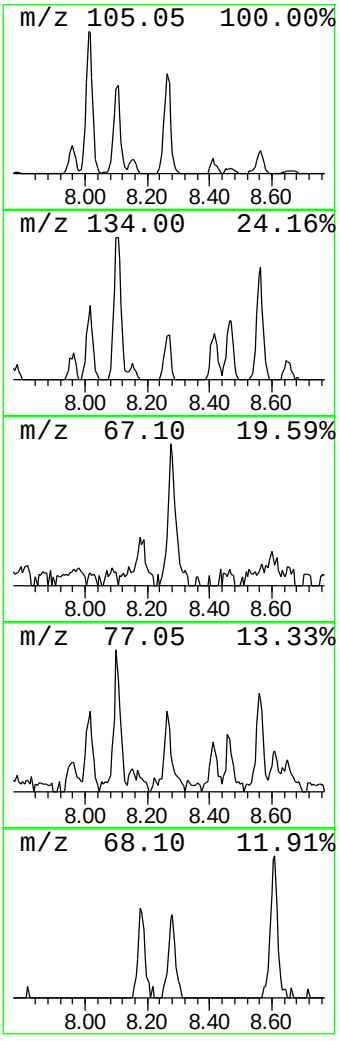
Peak Number 6 2-Tolyloxirane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.27	2.11 ng/ul	82629	1,4-Dichlorobenzene-d4	7.46

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Tolyloxirane	134	C9H10O	002783-26-8	64
2	Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	64
3	Benzene, (1-methylpropyl)-	134	C10H14	000135-98-8	60
4	Benzene, 1-methyl-2-propyl-	134	C10H14	001074-17-5	53
5	Benzeneacetaldehyde, .alpha.-met...	134	C9H10O	000093-53-8	53





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Acq On : 04 Jun 2020 19:12
Operator : JU/CG
Sample : L2844-15
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C0AF2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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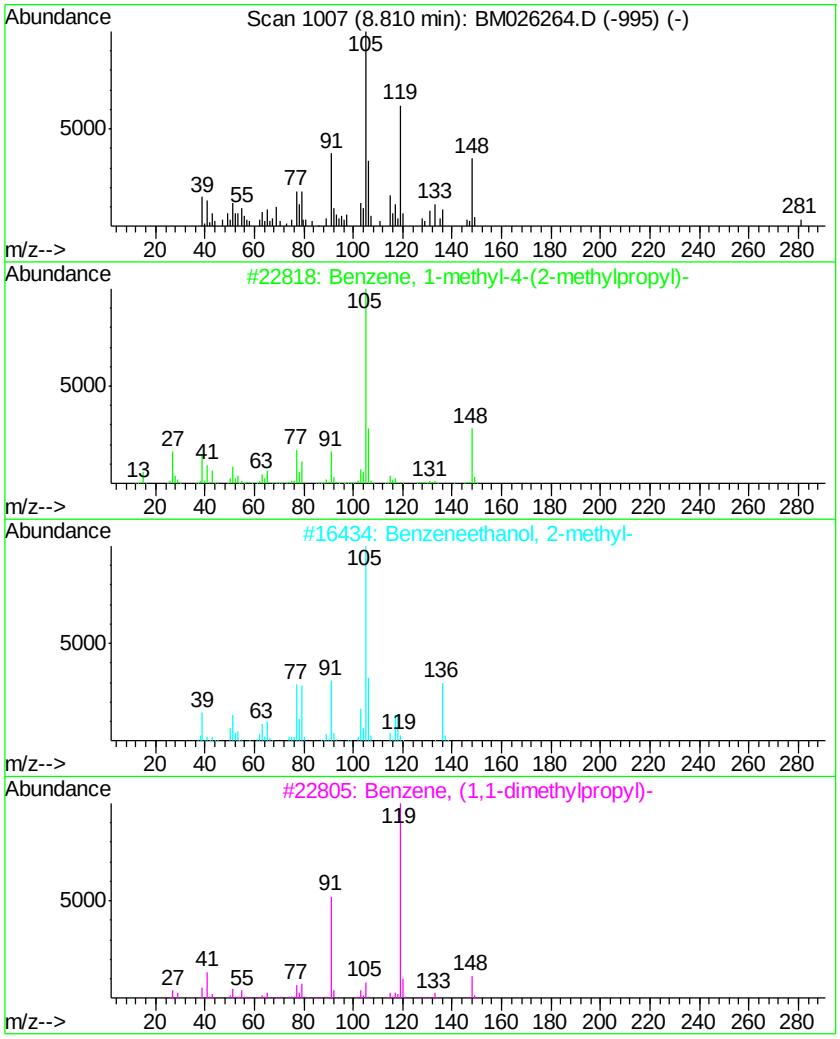
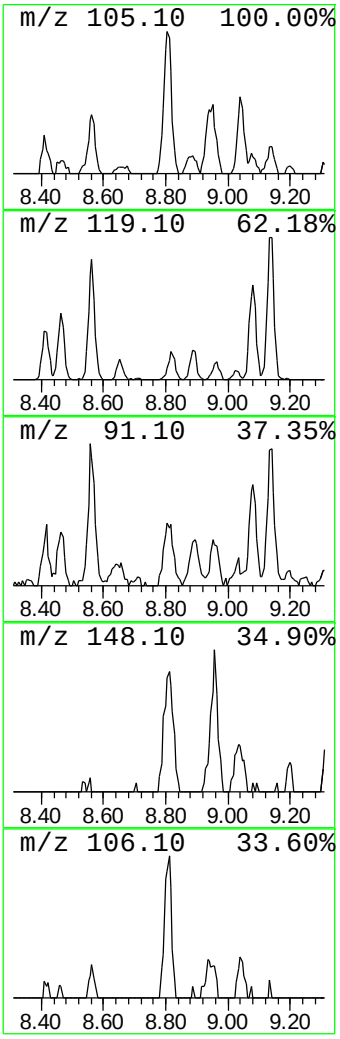
Peak Number 7 unknown-01 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.81	2.14 ng/ul	84003	1,4-Dichlorobenzene-d4	7.46

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-methyl-4-(2-methylpro...	148	C11H16	005161-04-6	46
2	Benzeneethanol, 2-methyl-	136	C9H12O	019819-98-8	41
3	Benzene, (1,1-dimethylpropyl)-	148	C11H16	002049-95-8	38
4	2-Butanone, 4-phenyl-	148	C10H12O	002550-26-7	38
5	3-Buten-2-ol, 4-phenyl-, (E)-	148	C10H12O	1000163-70-9	38





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Data File : BM026264.D
Acq On : 04 Jun 2020 19:12
Operator : JU/CG
Sample : L2844-15
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C0AF2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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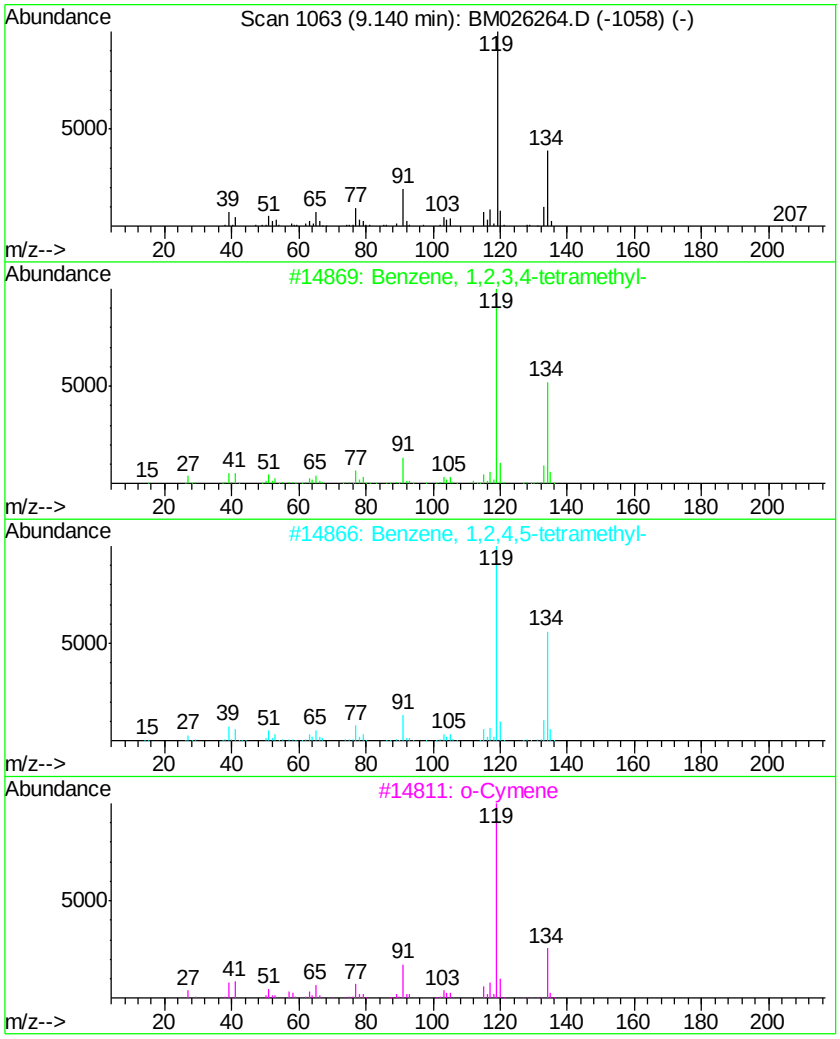
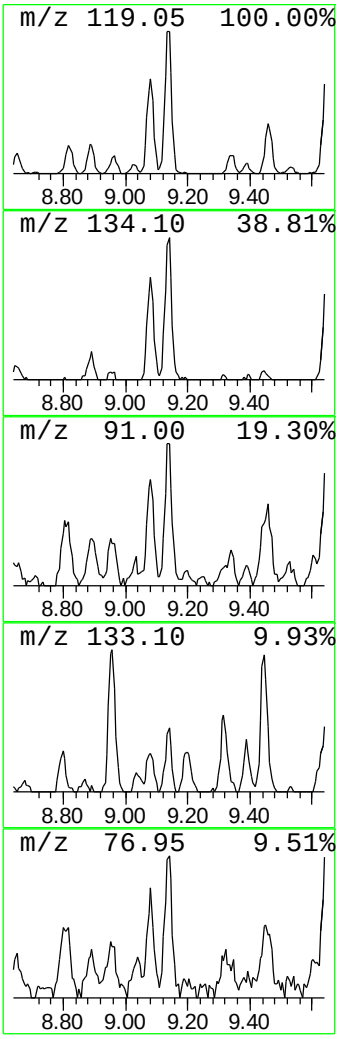
Peak Number 8 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.14	2.29 ng/ul	129041	Naphthalene-d8	10.22

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	94
2	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	94
3	o-Cymene	134	C10H14	000527-84-4	94
4	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	94
5	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	91





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Data File : BM026264.D
Acq On : 04 Jun 2020 19:12
Operator : JU/CG
Sample : L2844-15
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C0AF2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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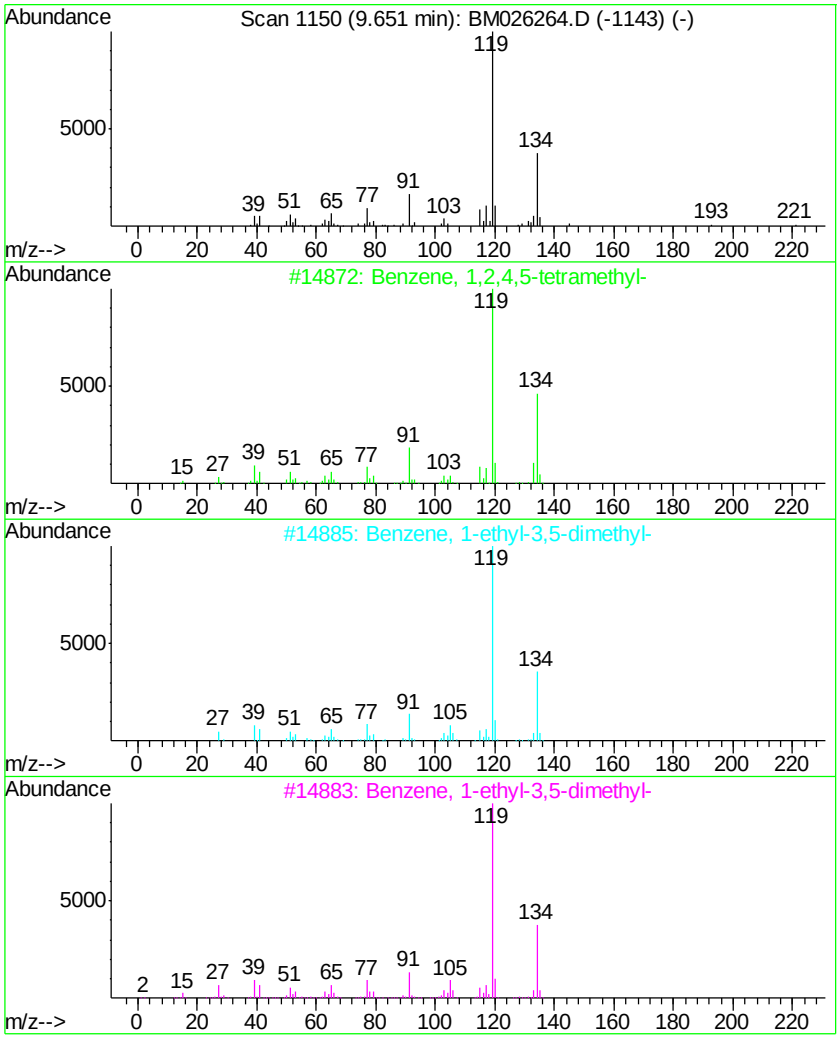
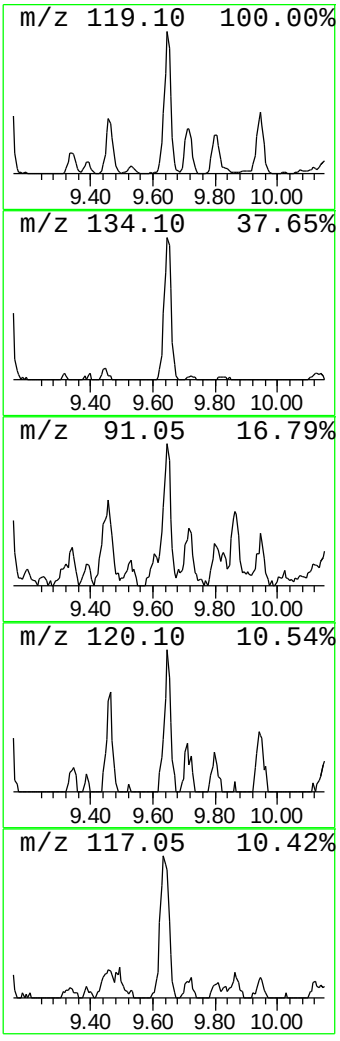
Peak Number 9 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.65	2.37 ng/ul	133542	Naphthalene-d8	10.22

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	93
2	Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	90
3	Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	90
4	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	90
5	o-Cymene	134	C10H14	000527-84-4	87





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM060420\
Data File : BM026264.D
Acq On : 04 Jun 2020 19:12
Operator : JU/CG
Sample : L2844-15
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C0AF2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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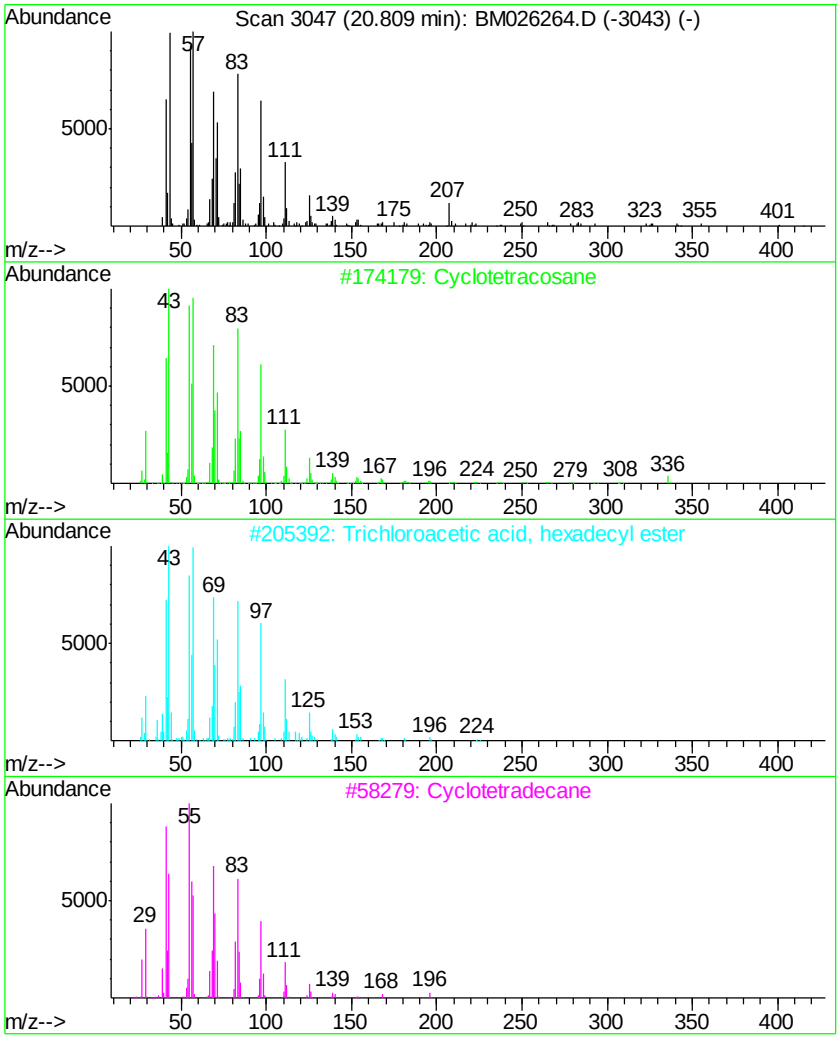
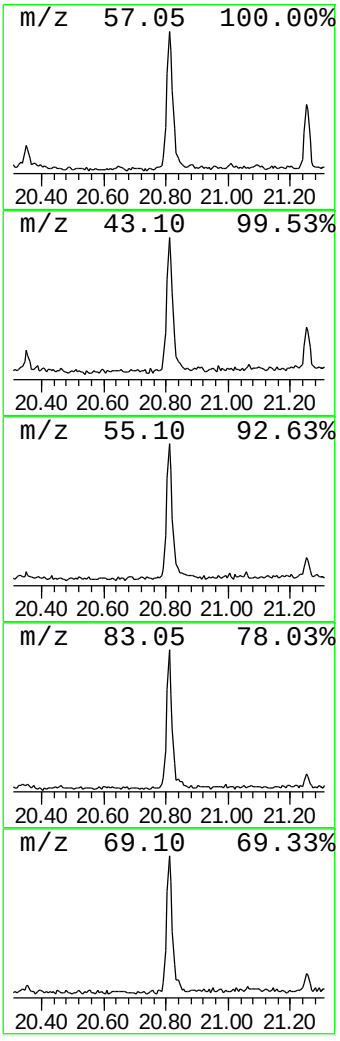
Peak Number 10 (DEL) Alkane: Cyclic20.81 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.81	2.72 ng/ul	248583	Chrysene-d12	21.06

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclotetracosane	336	C24H48	000297-03-0	95
2	Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	95
3	Cyclotetradecane	196	C14H28	000295-17-0	94
4	Heptafluorobutyric acid, pentade...	424	C19H31F7O2	959261-23-5	94
5	Cyclooctacosane	392	C28H56	000297-24-5	91





Tentatively Identified Compound (LSC) summary

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM060420\
Data File : BM026264.D
Acq On : 04 Jun 2020 19:12
Operator : JU/CG
Sample : L2844-15
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C0AF2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

					Internal Standard			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
(DEL) Alkane: Cyc...	5.78	2.1	ng/ul	83016	1	7.46	785069	20.0
Pentalene, octahy...	5.99	2.2	ng/ul	87358	1	7.46	785069	20.0
Benzene, 1,2,4-tr...	7.12	5.7	ng/ul	223742	1	7.46	785069	20.0
Benzene, 1,2,3-tr...	7.58	2.6	ng/ul	101436	1	7.46	785069	20.0
Benzene, 1-ethyl-...	8.10	3.3	ng/ul	130052	1	7.46	785069	20.0
2-Tolyloxirane	8.27	2.1	ng/ul	82629	1	7.46	785069	20.0
unknown-01	8.81	2.1	ng/ul	84003	1	7.46	785069	20.0
Benzene, 1,2,3,4-...	9.14	2.3	ng/ul	129041	2	10.22	1126960	20.0
Benzene, 1,2,4,5-...	9.65	2.4	ng/ul	133542	2	10.22	1126960	20.0
(DEL) Alkane: Cyc...	20.81	2.7	ng/ul	248583	5	21.06	1829100	20.0