LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062320\
Data File : BM026514.D
Acq On : 24 Jun 2020 00:29
Operator : JU/CG
Sample : L3069-05
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB7F6

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-BM061220MA.M
Title : SVOA CALIBRATION

Signal : TIC: BM026517.D

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.969	11	14	17	rBV	37829	45887	1.31%	0.113%
2	2.999	17	19	26	rVB	41431	55193	1.58%	0.136%
3	4.093	199	205	217	rVB	88030	144468	4.13%	0.356%
4	4.746	311	316	322	rBV	28478	46140	1.32%	0.114%
5	5.357	416	420	425	rVB4	7631	11211	0.32%	0.028%
6	5.704	469	479	483	rBV5	5424	14412	0.41%	0.036%
7	6.587	617	629	640	rBV	158862	298302	8.52%	0.735%
8	6.734	647	654	664	rBV	422836	660065	18.86%	1.626%
9	6.875	674	678	680	rBV3	6763	10361	0.30%	0.026%
10	6.922	680	686	694	rVV	503237	788853	22.54%	1.944%
11	6.987	694	697	704	rVV6	10502	18645	0.53%	0.046%
12	7.051	704	708	714	rVB4	10792	16091	0.46%	0.040%
13	7.240	732	740	746	rBV7	7959	20006	0.57%	0.049%
14	7.381	758	764	772	rVB	486892	733741	20.96%	1.808%
15	7.463	772	778	785	rBV4	22562	48813	1.39%	0.120%
16	7.698	812	818	822	rBV7	6229	12682	0.36%	0.031%
17	7.746	822	826	831	rVB2	22095	33690	0.96%	0.083%
18	8.046	867	877	880	rBV5	17778	35708	1.02%	0.088%
19	8.104	880	887	896	rVV	431952	690728	19.73%	1.702%
20	8.363	925	931	939	rBV	131259	209742	5.99%	0.517%
21	8.481	946	951	953	rBV	26944	38901	1.11%	0.096%
22	8.528	953	959	969	rVV	575851	955705	27.30%	2.355%
23	8.610	969	973	980	rVV8	9236	18753	0.54%	0.046%
24	8.722	983	992	996	rBV10	8697	22883	0.65%	0.056%
25	8.875	1011	1018	1023	rBV3	44170	74600	2.13%	0.184%
26	8.934	1024	1028	1033	rVB6	7281	12255	0.35%	0.030%
27	9.010	1033	1041	1049	rVB6	24040	49596	1.42%	0.122%
28	9.104	1051	1057	1062	rBV4	20370	32698	0.93%	0.081%
29	9.240	1074	1080	1090	rBV	520599	829111	23.69%	2.043%
30	9.363	1094	1101	1107	rBV6	15503	31895	0.91%	0.079%
31	9.463	1112	1118	1123	rVB7	15168	31091	0.89%	0.077%
32	9.669	1144	1153	1165	rVB9	28445	88424	2.53%	0.218%
33	9.775	1165	1171	1193	rBV	753511	1342719	38.36%	3.308%
34	9.951	1193	1201	1207	rBV10	13283	31743	0.91%	0.078%

LSC Area Percent Report

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Operator : JU/CG
Sample : L3069-05
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB7F6

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-BM061220MA.M
Title : SVOA CALIBRATION

35	10.134	1223	1232	1241	rBV	727724	1214634	34.70%	2.993%
36	10.287	1249	1258	1271	rBV	586239	1095464	31.30%	2.699%
37	10.404	1274	1278	1281	rVB3	12455	16530	0.47%	0.041%
38	10.492	1291	1293	1298	rVB6	8166	12008	0.34%	0.030%
39	10.551	1298	1303	1308	rVB8	6365	11568	0.33%	0.029%
40	10.604	1308	1312	1319	rVB10	6502	10656	0.30%	0.026%
41	10.722	1330	1332	1340	rVB9	10947	20260	0.58%	0.050%
42	10.887	1357	1360	1365	rBV2	15662	24791	0.71%	0.061%
43	10.981	1372	1376	1380	rVV6	10204	18226	0.52%	0.045%
44	11.345	1434	1438	1441	rBV7	14056	23324	0.67%	0.057%
45	11.439	1450	1454	1457	rBV3	9476	12881	0.37%	0.032%
46	11.510	1457	1466	1472	rVV	152266	259579	7.42%	0.640%
47	11.575	1474	1477	1479	rVB4	10650	13323	0.38%	0.033%
48	11.669	1486	1493	1498	rBV2	36656	65647	1.88%	0.162%
49	11.733	1500	1504	1507	rBV5	11582	17975	0.51%	0.044%
50	11.881	1523	1529	1538	rVB6	22441	50837	1.45%	0.125%
51	11.963	1538	1543	1548	rBV7	23694	45771	1.31%	0.113%
52	12.045	1554	1557	1563	rBV6	13463	26288	0.75%	0.065%
53	12.181	1576	1580	1586	rVB9	11984	24083	0.69%	0.059%
54	12.386	1609	1615	1620	rVB9	13781	28530	0.82%	0.070%
55	12.757	1672	1678	1686	rBV8	14639	45343	1.30%	0.112%
56	12.839	1688	1692	1695	rBV5	19002	28437	0.81%	0.070%
57	12.886	1697	1700	1707	rVB4	22245	43797	1.25%	0.108%
58	12.957	1707	1712	1714	rBV2	18933	30358	0.87%	0.075%
59	13.080	1728	1733	1741	rBV4	45510	75843	2.17%	0.187%
60	13.180	1746	1750	1753	rBV5	14900	23044	0.66%	0.057%
61	13.286	1765	1768	1775	rVB7	21131	35635	1.02%	0.088%
62	13.463	1793	1798	1803	rBV	1491173	1982647	56.64%	4.885%
63	13.510	1803	1806	1811	rVB	146531	183931	5.25%	0.453%
64	13.645	1825	1829	1836	rVB9	21574	35309	1.01%	0.087%
65	13.722	1836	1842	1849	rBV	1545386	2182028	62.34%	5.376%
66	13.880	1866	1869	1875	rVB2	28027	39702	1.13%	0.098%
67	13.963	1879	1883	1890	rVV3	66495	98647	2.82%	0.243%
68	14.033	1890	1895	1902	rVV	1102186	1657625	47.36%	4.084%
69	14.092	1903	1905	1909	rVB3	17617	22542	0.64%	0.056%
70	14.298	1935	1940	1952	rBV	107817	275760	7.88%	0.679%
71	14.657	1996	2001	2009	rBV	14589	42960	1.23%	0.106%

LSC Area Percent Report

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Operator : JU/CG
Sample : L3069-05
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB7F6

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-BM061220MA.M
Title : SVOA CALIBRATION

72	14.810	2022	2027	2043	rBV	769957	1081750	30.90%	2.665%
73	15.039	2060	2066	2074	rBV	1906990	2712716	77.50%	6.683%
74	15.186	2085	2091	2099	rBV	655493	979561	27.98%	2.413%
75	15.310	2109	2112	2117	rVB	95426	114250	3.26%	0.281%
76	15.574	2151	2157	2169	rVB	798259	1092008	31.20%	2.690%
77	15.751	2181	2187	2194	rBV3	78776	171063	4.89%	0.421%
78	15.833	2199	2201	2209	rVB8	30414	39360	1.12%	0.097%
79	15.910	2210	2214	2218	rVB2	42381	52323	1.49%	0.129%
80	16.086	2239	2244	2250	rBV	415048	563499	16.10%	1.388%
81	16.151	2252	2255	2263	rVB	79221	102190	2.92%	0.252%
82	16.369	2288	2292	2299	rVB	92485	131683	3.76%	0.324%
83	16.621	2332	2335	2340	rVB	40508	49626	1.42%	0.122%
84	16.757	2355	2358	2359	rBV3	29217	29842	0.85%	0.074%
85	16.792	2359	2364	2370	rVB	1456185	1956552	55.90%	4.820%
86	16.892	2375	2381	2391	rBV	2151223	2962476	84.63%	7.299%
87	17.733	2521	2524	2528	rBV4	36236	50319	1.44%	0.124%
88	18.127	2588	2591	2595	rBV2	28379	37102	1.06%	0.091%
89	18.486	2649	2652	2658	rVB3	64651	93687	2.68%	0.231%
90	18.833	2708	2711	2715	rVB2	27662	32781	0.94%	0.081%
91	19.198	2768	2773	2780	rBV	2565371	3500401	100.00%	8.624%
92	19.274	2781	2786	2792	rVV	474380	590347	16.87%	1.454%
93	20.280	2955	2957	2962	rVB2	71607	73609	2.10%	0.181%
94	20.756	3034	3038	3046	rBV	419789	509784	14.56%	1.256%
95	21.003	3075	3080	3090	rBV	1808222	2128178	60.80%	5.243%
96	22.509	3333	3336	3341	rVB	107286	131900	3.77%	0.325%
97	22.962	3407	3413	3419	rBV	1186056	1961518	56.04%	4.833%
98	23.021	3421	3423	3428	rVV	150292	199485	5.70%	0.491%
99	23.092	3429	3435	3444	rVB	1044267	1766101	50.45%	4.351%
100	23.603	3519	3522	3529	rVB	135286	223084	6.37%	0.550%

Sum of corrected areas: 40588290

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Data File : BM026514.D
Acq On : 24 Jun 2020 00:29
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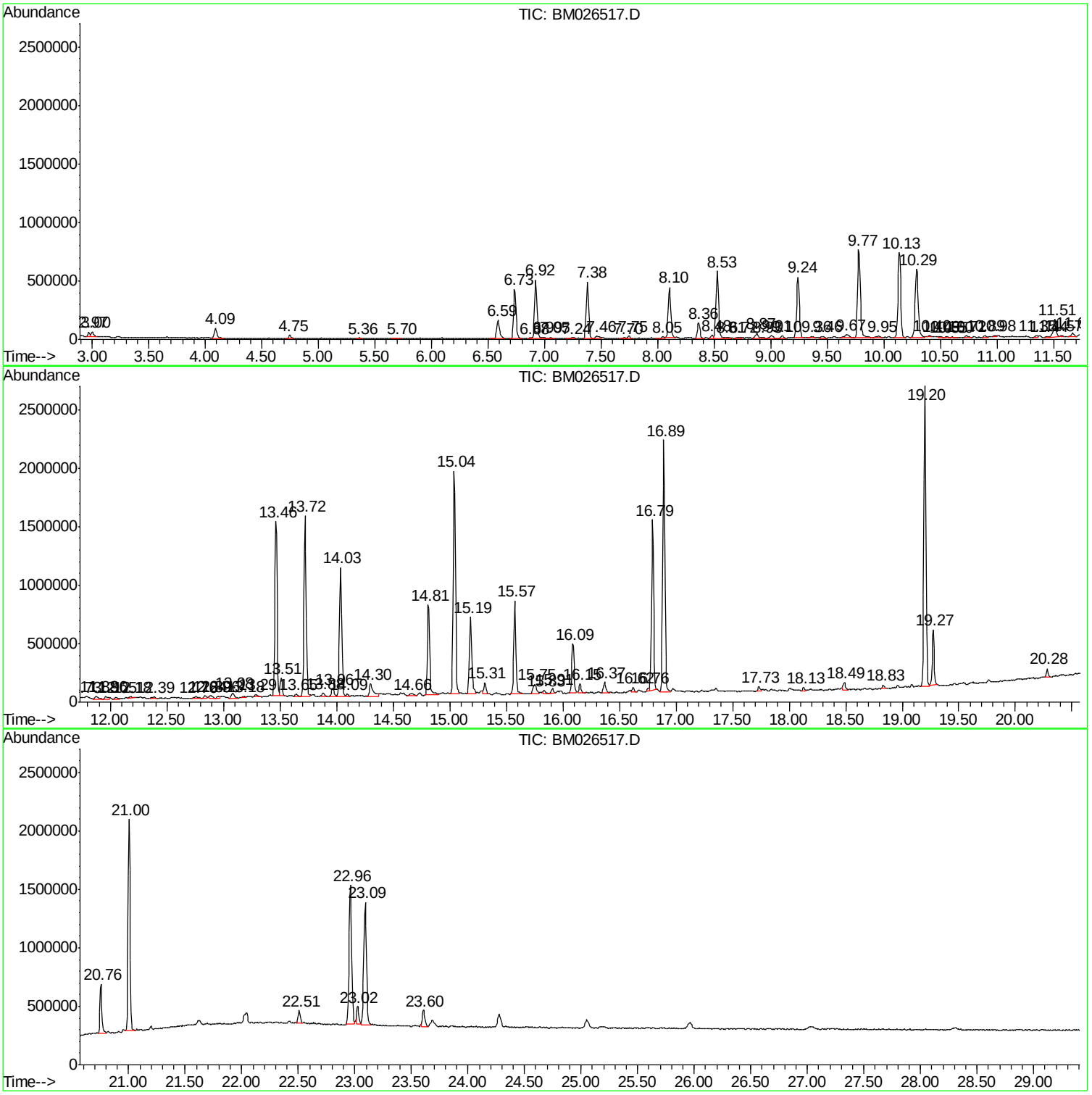
Sample : L3069-05
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
CB7F6

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061220MA.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\BM062320\
Data File : BM026514.D
Acq On : 24 Jun 2020 00:29
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Instrument : BNA_M
ClientSampleId : CB7F6

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061220MA.M
Quant Title : SVOA CALIBRATION

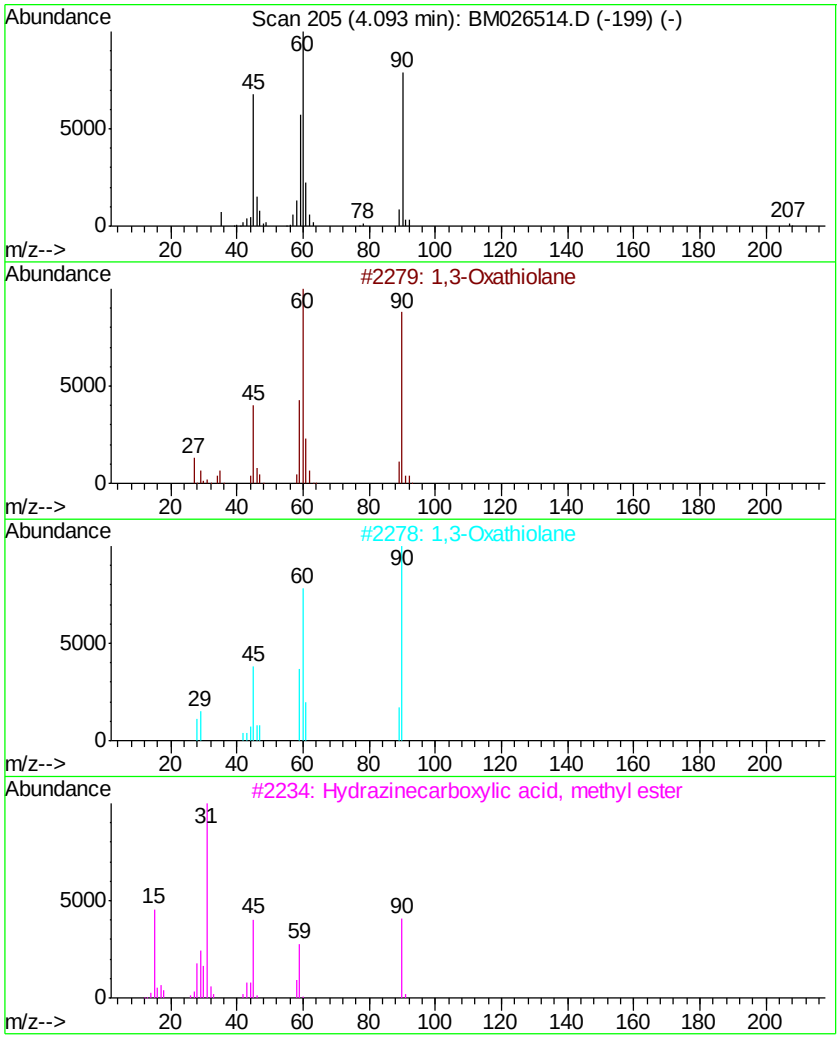
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

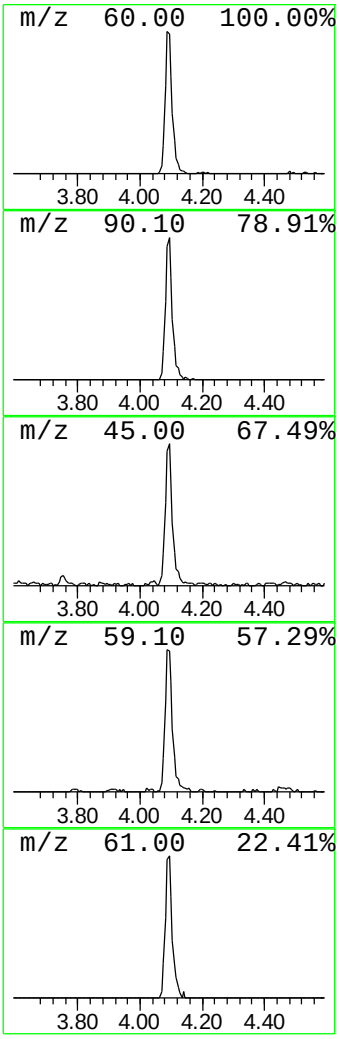
****

Peak Number 1 1,3-Oxathiolane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.09	3.94 ng/ul	144468	1,4-Dichlorobenzene-d4	7.38

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,3-Oxathiolane	90	C3H6OS	002094-97-5	87
2	1,3-Oxathiolane	90	C3H6OS	002094-97-5	64
3	Hydrazinecarboxylic acid, methyl...	90	C2H6N2O2	006294-89-9	59
4	3-Thietanol	90	C3H6OS	010304-16-2	53
5	Thiourea, methyl-	90	C2H6N2S	000598-52-7	50





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062320\
Data File : BM026514.D
Acq On : 24 Jun 2020 00:29
Operator : JU/CG
Sample : L3069-05
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB7F6

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061220MA.M
Quant Title : SVOA CALIBRATION

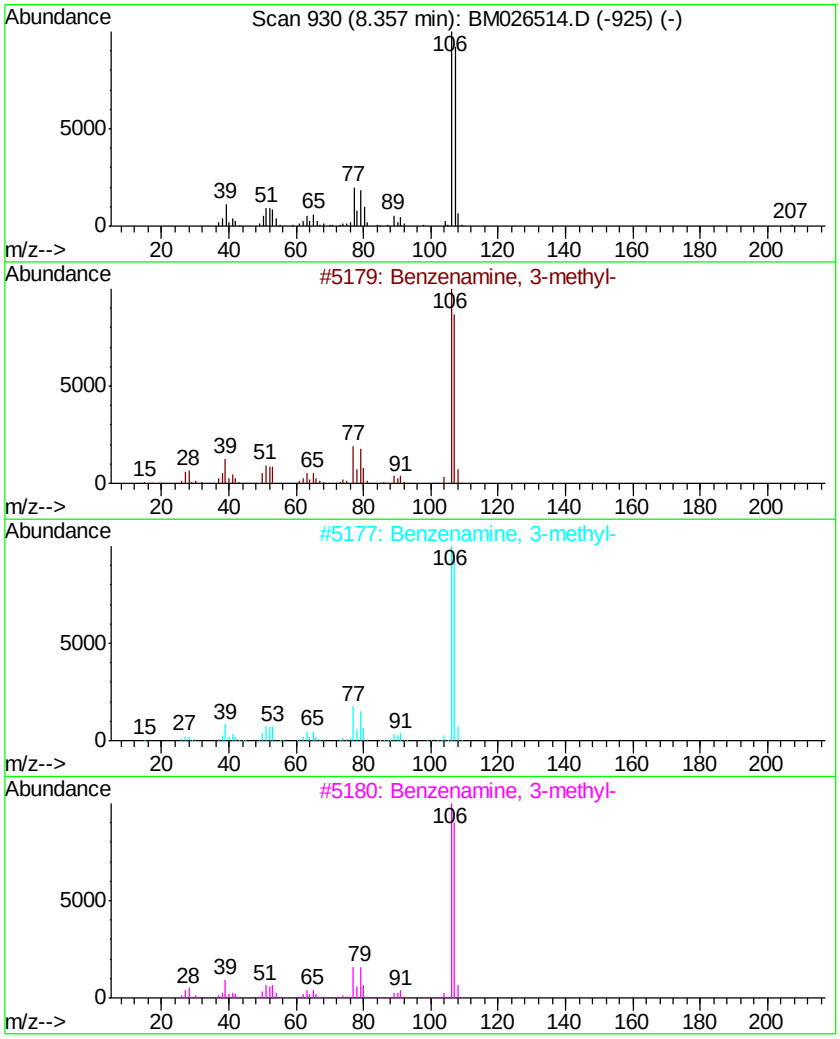
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

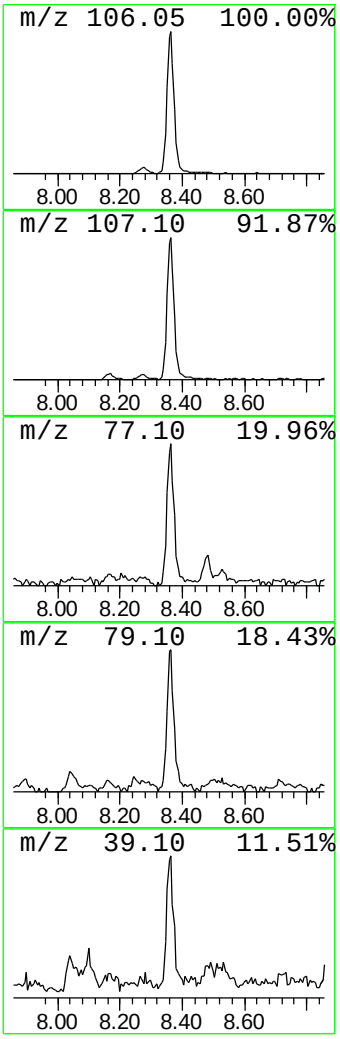
****

Peak Number 2 Benzenamine, 3-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.36	5.72 ng/ul	209742	1,4-Dichlorobenzene-d4	7.38

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzenamine, 3-methyl-	107	C7H9N	000108-44-1	96
2	Benzenamine, 3-methyl-	107	C7H9N	000108-44-1	95
3	Benzenamine, 3-methyl-	107	C7H9N	000108-44-1	95
4	Benzenamine, 3-methyl-	107	C7H9N	000108-44-1	93
5	Benzenamine, 3-methyl-	107	C7H9N	000108-44-1	93





Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062320\
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Operator : JU/CG
Sample : L3069-05
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB7F6

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061220MA.M
Quant Title : SVOA CALIBRATION

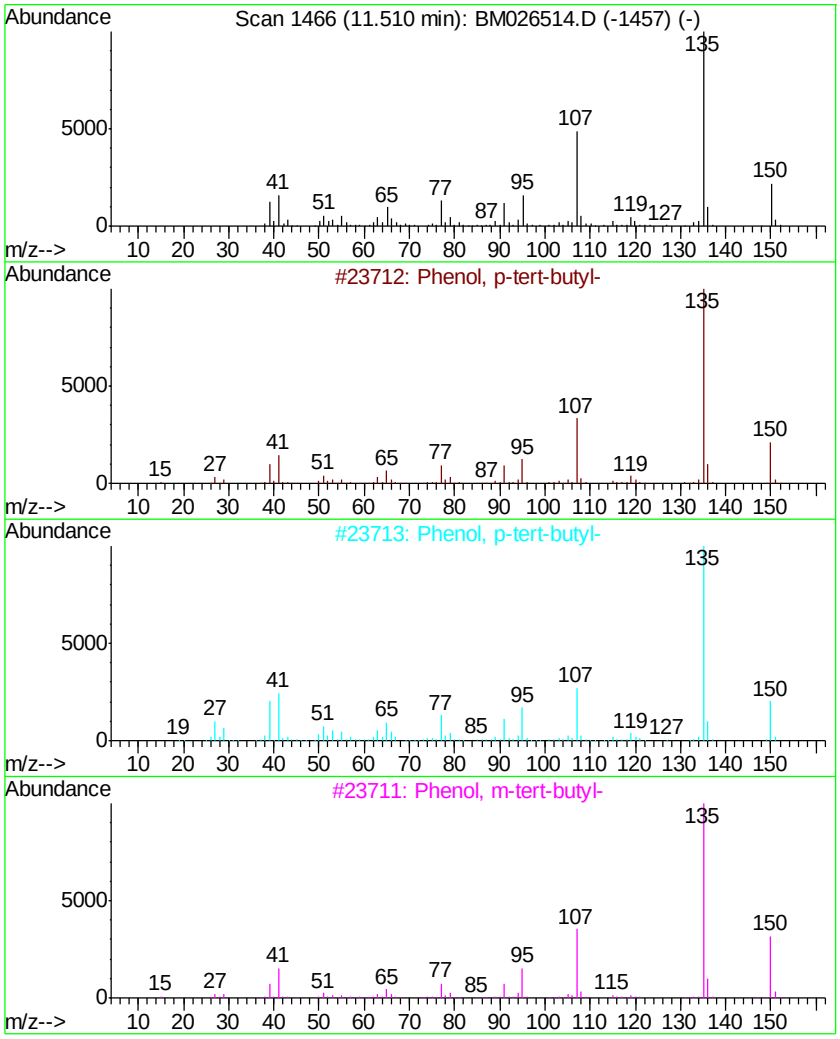
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

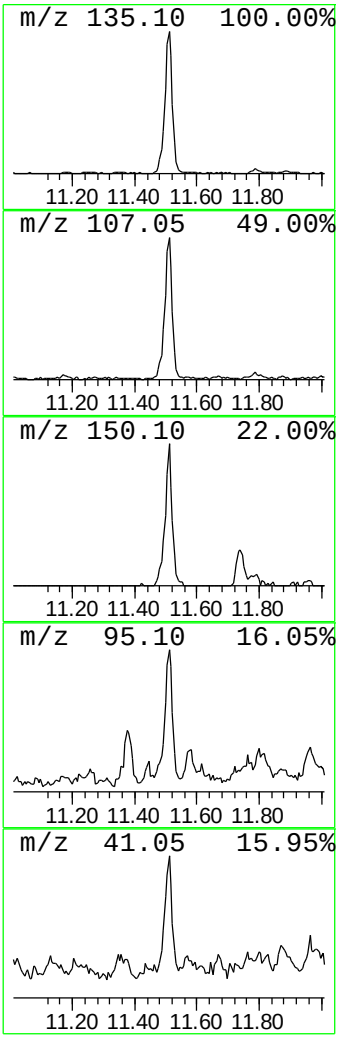
****

Peak Number 3 Phenol, p-tert-butyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.51	4.27 ng/ul	259579	Naphthalene-d8	10.13

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	96
2	Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	94
3	Phenol, m-tert-butyl-	150	C10H14O	000585-34-2	94
4	Phenol, m-tert-butyl-	150	C10H14O	000585-34-2	93
5	Phenol, 2-(1,1-dimethylethyl)-	150	C10H14O	000088-18-6	90





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062320\
Data File : BM026514.D
Acq On : 24 Jun 2020 00:29
Operator : JU/CG
Sample : L3069-05
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB7F6

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061220MA.M
Quant Title : SVOA CALIBRATION

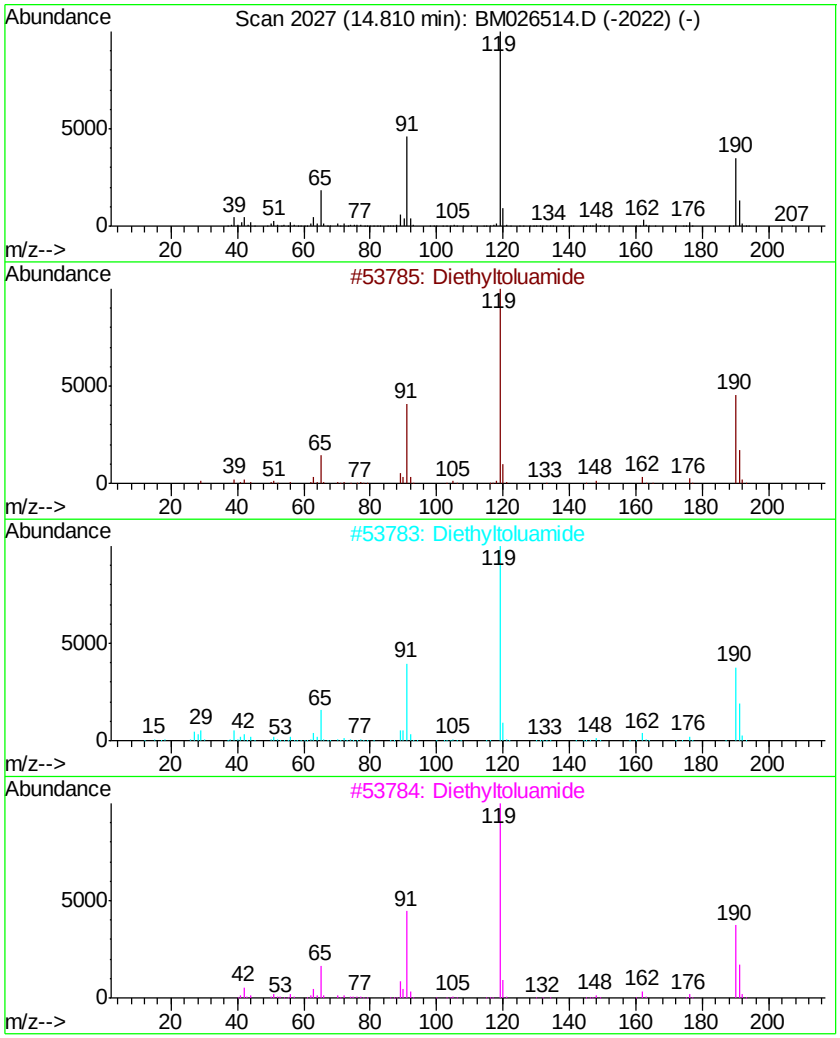
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

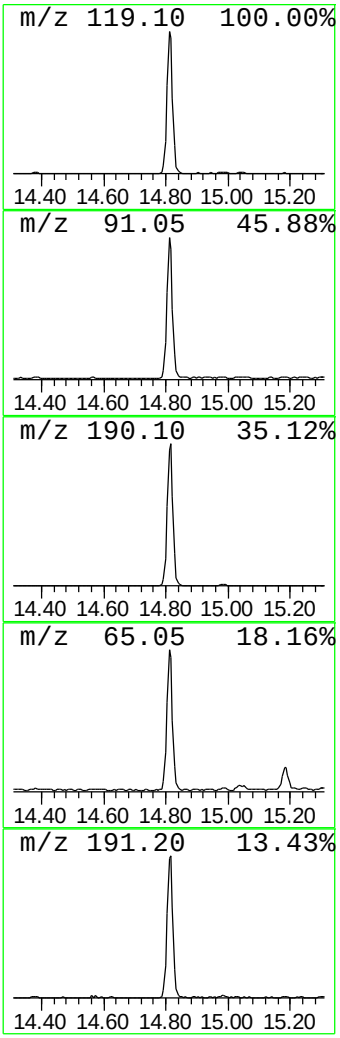
****

Peak Number 4 Diethyltoluamide Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.81	13.05 ng/ul	1081750	Acenaphthene-d10	14.03

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Diethyltoluamide	191	C12H17NO	000134-62-3	97
2	Diethyltoluamide	191	C12H17NO	000134-62-3	95
3	Diethyltoluamide	191	C12H17NO	000134-62-3	92
4	1-Propanone, 2-methyl-1-(2-methy...	162	C11H14O	002040-21-3	72
5	L-Valine, N-(4-methylbenzoyl)-, ...	277	C16H23NO3	1000346-63-8	72





Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062320\
Data File : BM026514.D
Acq On : 24 Jun 2020 00:29
Operator : JU/CG
Sample : L3069-05
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB7F6

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061220MA.M
Quant Title : SVOA CALIBRATION

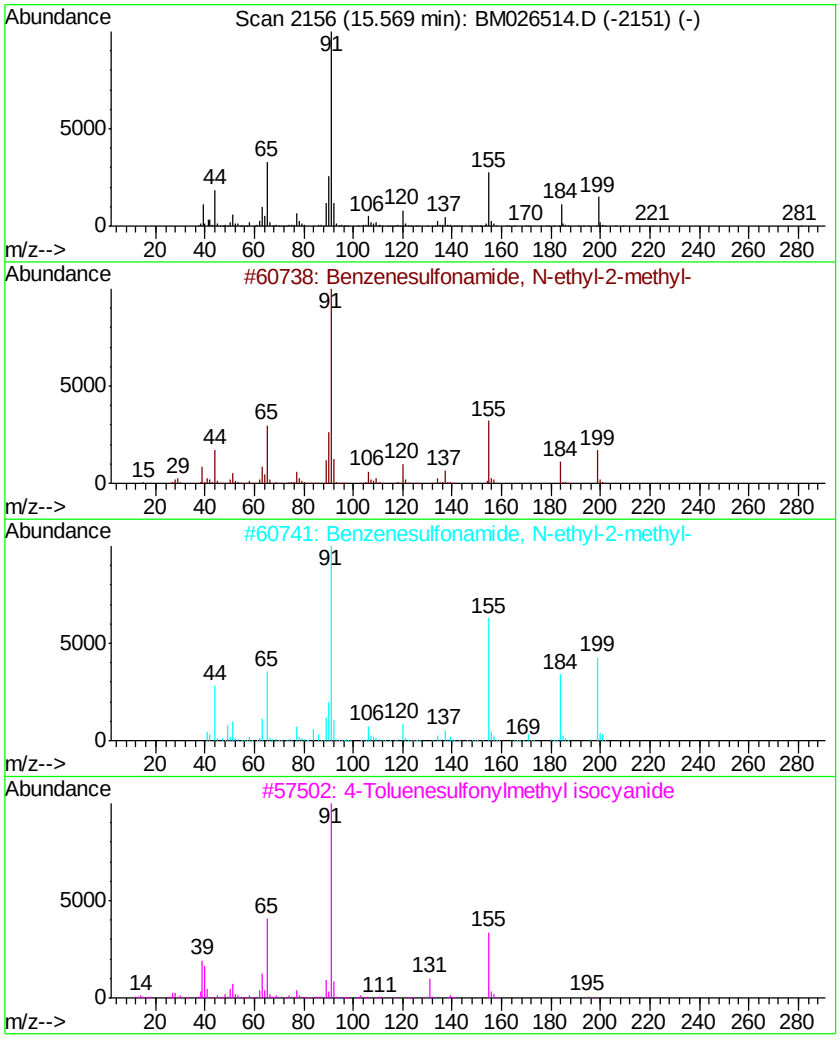
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

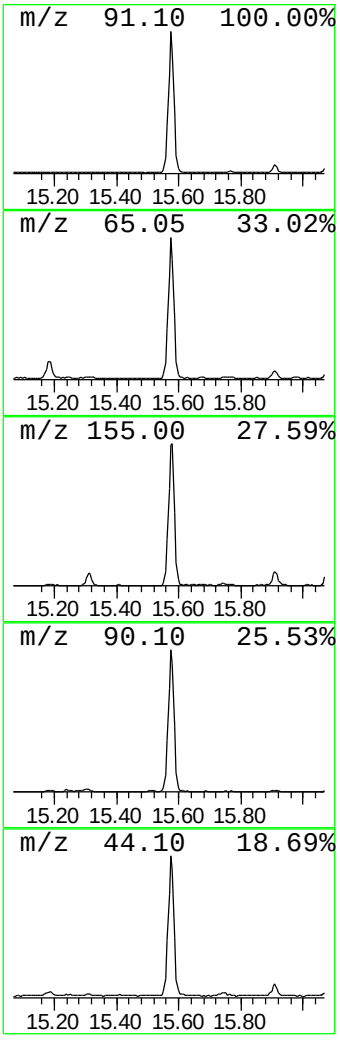
****

Peak Number 5 Benzenesulfonamide, N-ethyl... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.57	11.16 ng/ul	1092010	Phenanthrene-d10	16.79

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzenesulfonamide, N-ethyl-2-me...	199	C9H13NO2S	001077-56-1	98
2	Benzenesulfonamide, N-ethyl-2-me...	199	C9H13NO2S	001077-56-1	64
3	4-Toluenesulfonylmethyl isocyanide	195	C9H9NO2S	036635-61-7	58
4	Tolbutamide	270	C12H18N2O3S	000064-77-7	50
5	p-Toluenesulfonylacetonitrile	195	C9H9NO2S	005697-44-9	50





Library Search Compound Report

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Data File : BM026514.D
Acq On : 24 Jun 2020 00:29
Operator : JU/CG
Sample : L3069-05
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB7F6

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061220MA.M
Quant Title : SVOA CALIBRATION

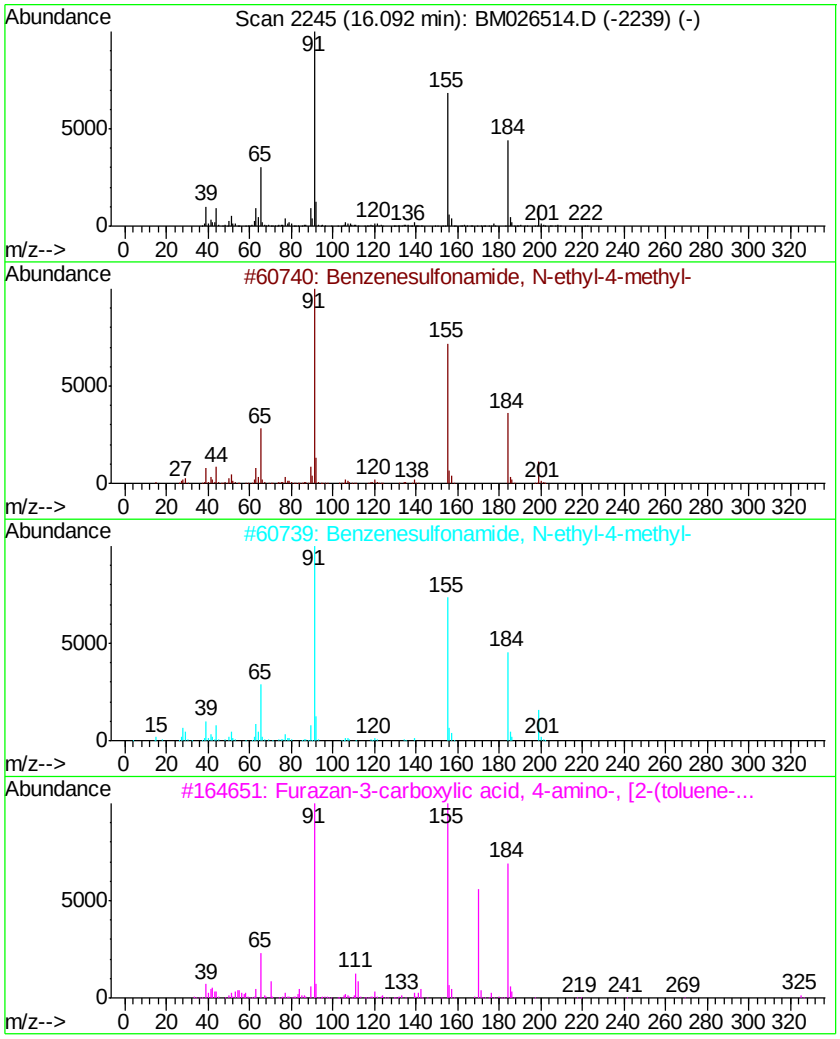
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

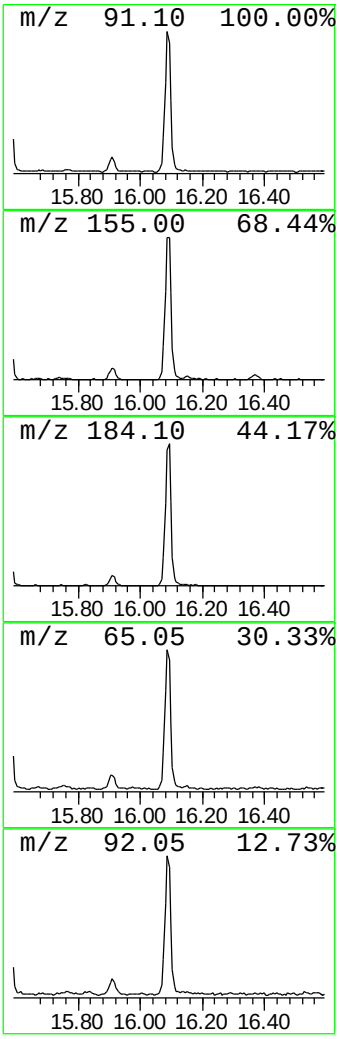
****

Peak Number 6 Benzenesulfonamide, N-ethyl... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.09	5.76 ng/ul	563499	Phenanthrene-d10	16.79

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzenesulfonamide, N-ethyl-4-me...	199	C9H13NO2S	000080-39-7	98
2	Benzenesulfonamide, N-ethyl-4-me...	199	C9H13NO2S	000080-39-7	96
3	Furazan-3-carboxylic acid, 4-ami...	325	C12H15N5O4S	1000304-69-4	78
4	4-Methyl-N-(2-oxobutyl)benzenesu...	241	C11H15NO3S	1000192-14-5	78
5	(O-p-Tosylisonitroso)malononitrile	249	C10H7N3O3S	020893-01-0	47





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Data File : BM026514.D
Acq On : 24 Jun 2020 00:29
Operator : JU/CG
Sample : L3069-05
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB7F6

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061220MA.M
Quant Title : SVOA CALIBRATION

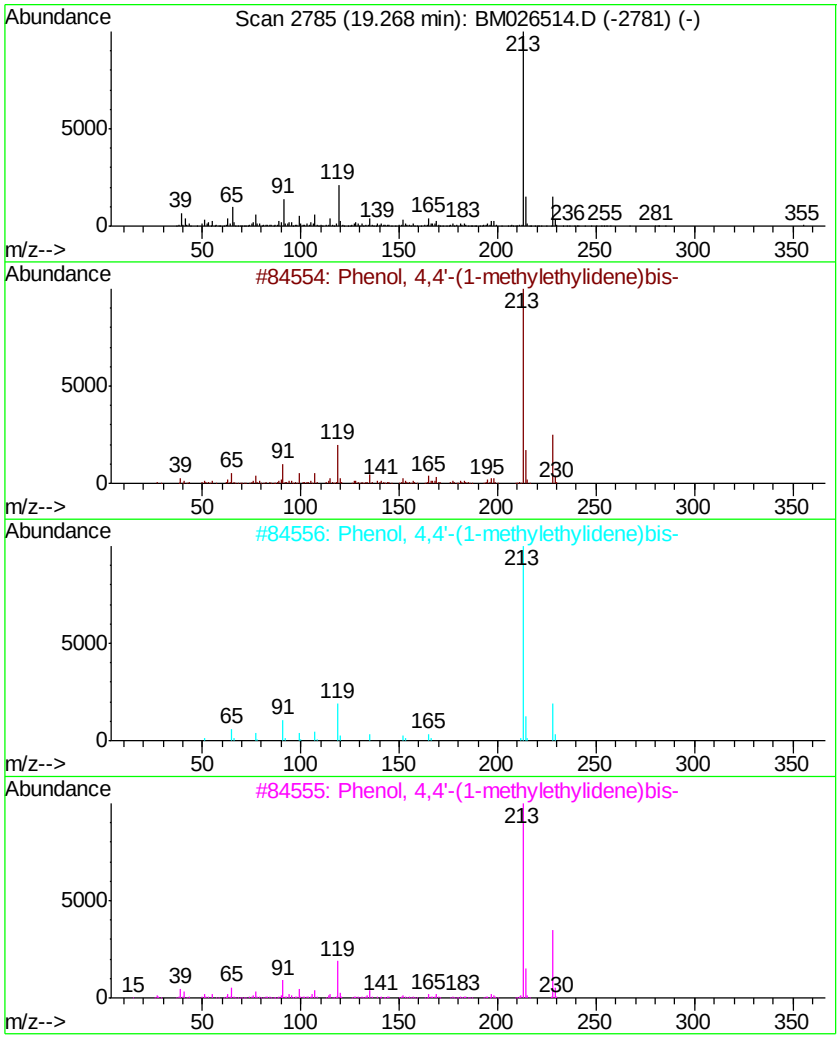
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

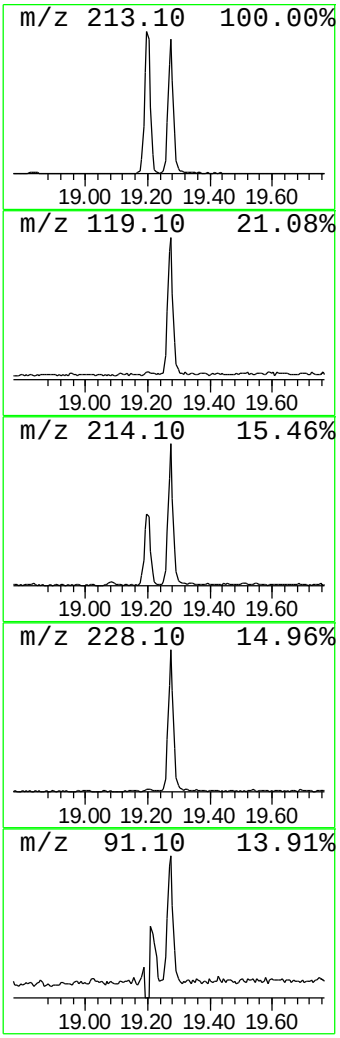
****

Peak Number 7 Phenol, 4,4'-(1-methylethyl... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.27	5.55 ng/ul	590347	Chrysene-d12	21.00

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 4,4'-(1-methylethylidene...	228	C15H16O2	000080-05-7	97
2	Phenol, 4,4'-(1-methylethylidene...	228	C15H16O2	000080-05-7	97
3	Phenol, 4,4'-(1-methylethylidene...	228	C15H16O2	000080-05-7	86
4	2H,8H-Benzo[1,2-b:5,4-b']dipyran...	228	C14H12O3	000553-19-5	64
5	2H,8H-Benzo[1,2-b:3,4-b']dipyran...	228	C14H12O3	000523-59-1	59





Library Search Compound Report

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Data File : BM026514.D
Acq On : 24 Jun 2020 00:29
Operator : JU/CG
Sample : L3069-05
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB7F6

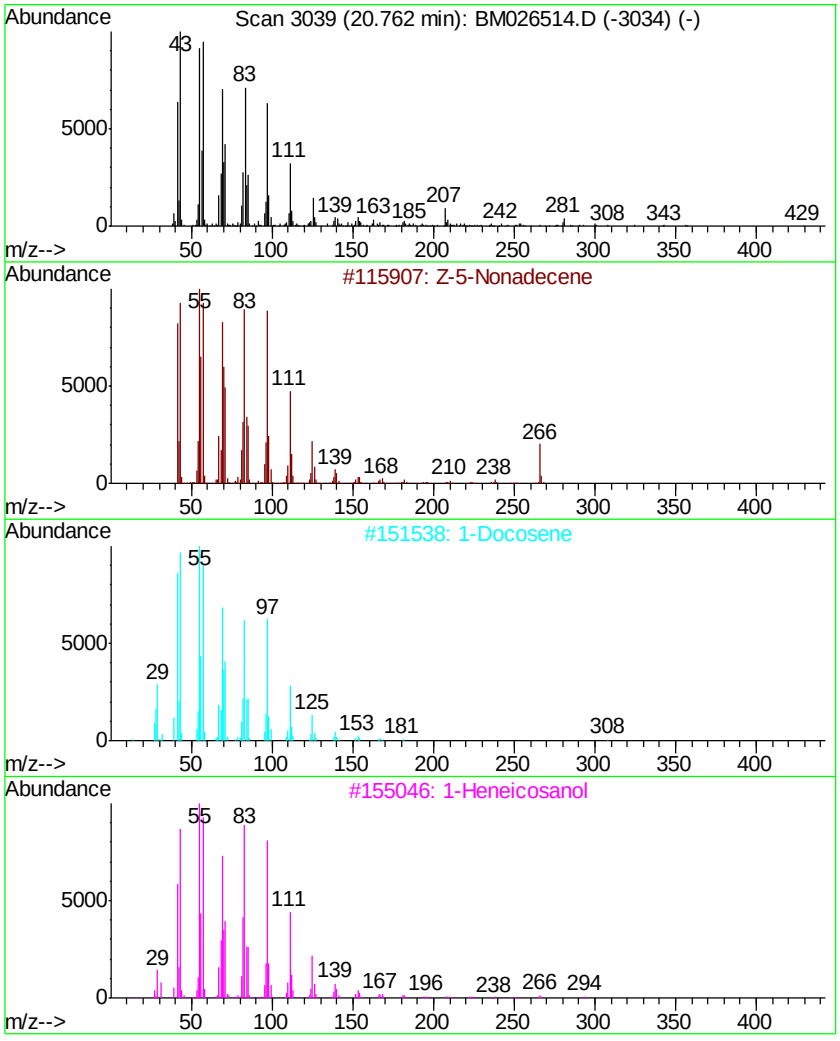
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061220MA.M
Quant Title : SVOA CALIBRATION

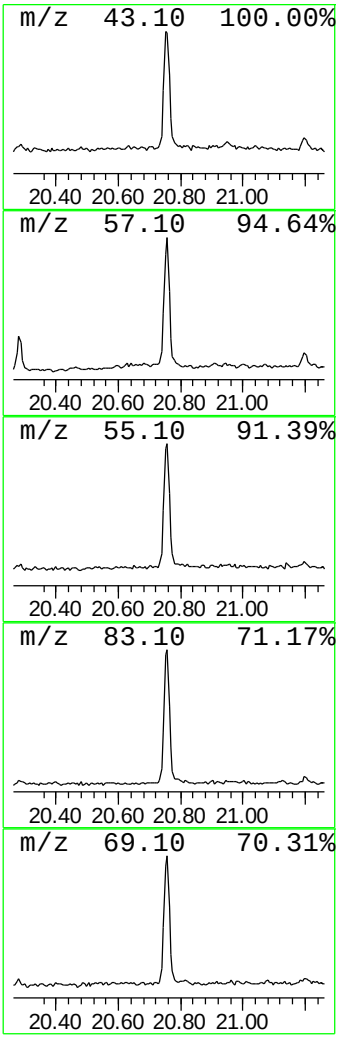
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.76	4.79 ng/ul	509784	Chrysene-d12	21.00

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Z-5-Nonadecene	266	C19H38	1000131-11-8	98
2	1-Docosene	308	C22H44	001599-67-3	95
3	1-Heneicosanol	312	C21H44O	015594-90-8	94
4	Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	94
5	Heptafluorobutyric acid, pentade...	424	C19H31F7O2	959261-23-5	94





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Data File : BM026514.D
Acq On : 24 Jun 2020 00:29
Operator : JU/CG
Sample : L3069-05
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB7F6

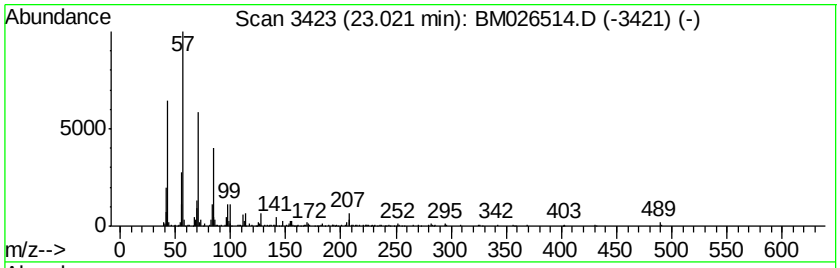
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061220MA.M
Quant Title : SVOA CALIBRATION

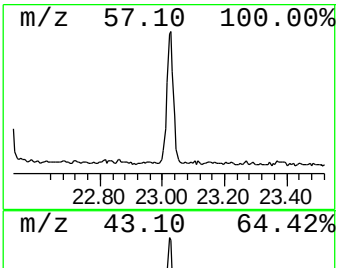
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.02	2.26 ng/ul	199485	Perylene-d12	23.09

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octacosane	394	C28H58	000630-02-4	90
2	Heneicosane, 11-decyl-	437	C31H64	055320-06-4	87
3	Tetratetracontane	619	C44H90	007098-22-8	87
4	Eicosane, 7-hexyl-	366	C26H54	055333-99-8	87
5	Tetracosane	338	C24H50	000646-31-1	86





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062320\
Data File : BM026514.D
Acq On : 24 Jun 2020 00:29
Operator : JU/CG
Sample : L3069-05
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB7F6

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061220MA.M
Quant Title : SVOA CALIBRATION

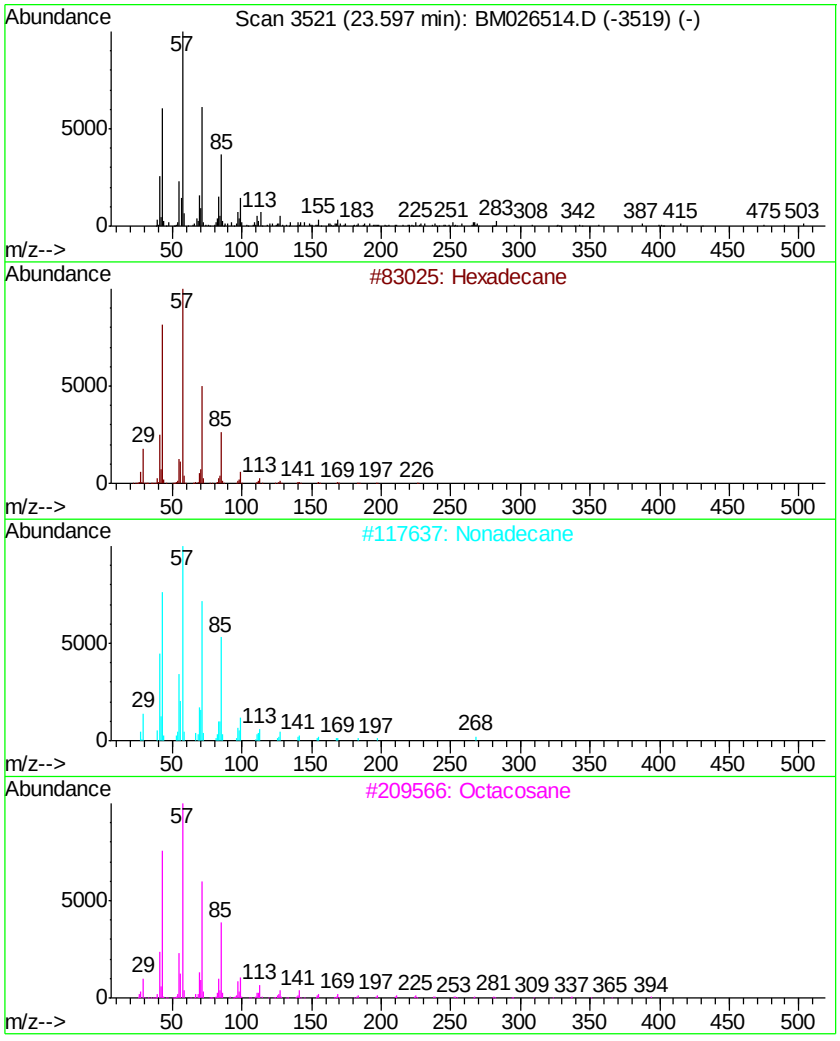
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

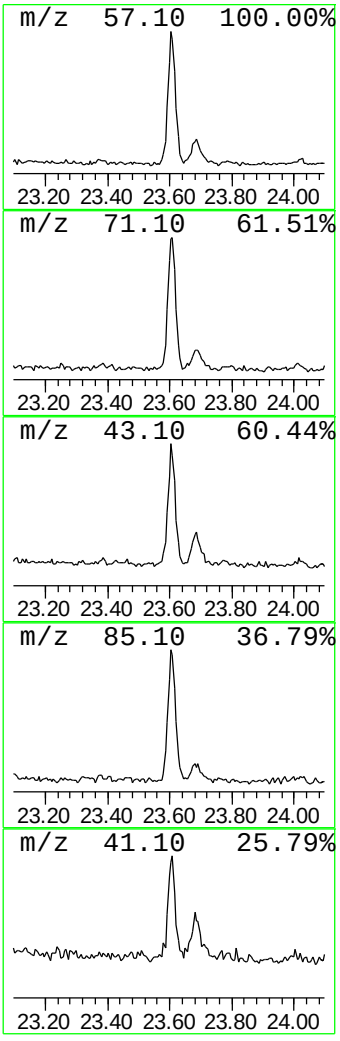
****

Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.60	2.53 ng/ul	223084	Perylene-d12	23.09

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane	226	C16H34	000544-76-3	93
2	Nonadecane	268	C19H40	000629-92-5	93
3	Octacosane	394	C28H58	000630-02-4	87
4	Tetracosane	338	C24H50	000646-31-1	87
5	Heneicosane	296	C21H44	000629-94-7	87





Tentatively Identified Compound (LSC) summary

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM062320\
Data File : BM026514.D
Acq On : 24 Jun 2020 00:29
Operator : JU/CG
Sample : L3069-05
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB7F6

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061220MA.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
1,3-Oxathiolane	4.09	3.9	ng/ul	144468	1	7.38	733741	20.0
Benzenamine, 3-me...	8.36	5.7	ng/ul	209742	1	7.38	733741	20.0
Phenol, p-tert-bu...	11.51	4.3	ng/ul	259579	2	10.13	1214630	20.0
Diethyltoluamide	14.81	13.1	ng/ul	1081750	3	14.03	1657630	20.0
Benzenesulfonamid...	15.57	11.2	ng/ul	1092010	4	16.79	1956550	20.0
Benzenesulfonamid...	16.09	5.8	ng/ul	563499	4	16.79	1956550	20.0
Phenol, 4,4'-(1-m...	19.27	5.5	ng/ul	590347	5	21.00	2128180	20.0
(DEL) Alkane: Str...	20.76	4.8	ng/ul	509784	5	21.00	2128180	20.0
(DEL) Alkane: Str...	23.02	2.3	ng/ul	199485	6	23.09	1766100	20.0
(DEL) Alkane: Str...	23.60	2.5	ng/ul	223084	6	23.09	1766100	20.0