LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG101420\
Data File : BG046896.D
Acq On : 14 Oct 2020 17:06
Operator : CG/JU
Sample : L4359-07
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : CB7N5

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_G\METHODS\SOM-EPA-BG092920MA.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	3.517	25	30	33	rBV2	14248	19560	1.40%	0.114%
2	3.552	33	36	44	rVB2	18706	28484	2.04%	0.166%
3	3.793	75	77	80	rVB4	3662	3831	0.27%	0.022%
4	4.110	127	131	135	rVB4	2348	3738	0.27%	0.022%
5	4.275	154	159	165	rVB4	7886	12853	0.92%	0.075%
6	4.386	175	178	181	rVB4	3257	3702	0.26%	0.022%
7	4.568	203	209	210	rBV3	2129	2690	0.19%	0.016%
8	4.715	228	234	239	rBV	42332	69032	4.93%	0.403%
9	5.185	311	314	316	rBV2	2772	2513	0.18%	0.015%
10	5.403	345	351	359	rVB	19875	31388	2.24%	0.183%
11	5.667	391	396	398	rBV3	2781	4595	0.33%	0.027%
12	5.802	416	419	422	rBV	1905	2084	0.15%	0.012%
13	5.926	436	440	444	rBV2	1938	3078	0.22%	0.018%
14	6.026	453	457	463	rVB4	5855	7862	0.56%	0.046%
15	6.355	509	513	519	rVB5	4511	8632	0.62%	0.050%
16	6.590	547	553	555	rBV3	4603	6018	0.43%	0.035%
17	6.643	559	562	565	rBV3	2185	3239	0.23%	0.019%
18	6.684	565	569	573	rVB3	3174	5998	0.43%	0.035%
19	6.731	573	577	578	rBV2	2997	3318	0.24%	0.019%
20	7.165	645	651	655	rVB3	1282	3037	0.22%	0.018%
21	7.242	658	664	674	rBV3	48707	105943	7.57%	0.618%
22	7.418	686	694	701	rBV	117347	193306	13.81%	1.128%
23	7.530	711	713	716	rVB3	3001	2522	0.18%	0.015%
24	7.571	716	720	721	rBV2	1913	2762	0.20%	0.016%
25	7.624	721	729	736	rVV	142572	247882	17.71%	1.446%
26	7.694	738	741	746	rVB6	4711	7726	0.55%	0.045%
27	7.747	746	750	751	rBV3	4269	3678	0.26%	0.021%
28	7.953	781	785	787	rBV3	3038	2673	0.19%	0.016%
29	7.988	790	791	798	rVB5	3431	3673	0.26%	0.021%
30	8.053	798	802	803	rBV	2365	2192	0.16%	0.013%
31	8.100	803	810	816	rVV	217567	369796	26.42%	2.158%
32	8.164	816	821	829	rVB5	11212	23497	1.68%	0.137%
33	8.235	829	833	837	rBV3	4032	6272	0.45%	0.037%
34	8.476	870	874	880	rVB	12562	18038	1.29%	0.105%

LSC Area Percent Report

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Sample : L4359-07
Misc :
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Instrument : BNA_G
ClientSampleId : CB7N5

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_G\METHODS\SOM-EPA-BG092920MA.M
Title : SVOA CALIBRATION

35	8.552	880	887	889	rBV3	3888	6544	0.47%	0.038%
36	8.634	900	901	905	rVB4	2665	2841	0.20%	0.017%
37	8.740	914	919	921	rBV5	11325	13905	0.99%	0.081%
38	8.787	922	927	937	rVB	120782	225017	16.08%	1.313%
39	8.887	941	944	948	rBV3	3006	4877	0.35%	0.028%
40	8.987	957	961	962	rBV3	4325	5252	0.38%	0.031%
41	9.075	969	976	985	rVB	106749	188890	13.50%	1.102%
42	9.192	992	996	1000	rBV3	15230	23785	1.70%	0.139%
43	9.257	1000	1007	1016	rVB	149967	273234	19.52%	1.594%
44	9.328	1016	1019	1020	rBV3	3316	3380	0.24%	0.020%
45	9.551	1056	1057	1060	rVB3	2904	2123	0.15%	0.012%
46	9.610	1062	1067	1074	rBV5	19996	39791	2.84%	0.232%
47	9.686	1075	1080	1086	rVB6	13459	23963	1.71%	0.140%
48	9.851	1103	1108	1113	rBV5	8984	15923	1.14%	0.093%
49	9.927	1119	1121	1124	rVB2	3610	3408	0.24%	0.020%
50	9.980	1124	1130	1141	rBV	142616	250876	17.92%	1.464%
51	10.109	1146	1152	1157	rBV7	9262	21437	1.53%	0.125%
52	10.168	1160	1162	1164	rBV2	3983	4597	0.33%	0.027%
53	10.274	1174	1180	1183	rBV3	6982	11022	0.79%	0.064%
54	10.321	1183	1188	1191	rVV6	8667	14337	1.02%	0.084%
55	10.385	1191	1199	1204	rVV6	6889	15602	1.11%	0.091%
56	10.438	1204	1208	1215	rVB8	17835	41121	2.94%	0.240%
57	10.520	1215	1222	1235	rBV	236690	448461	32.04%	2.617%
58	10.673	1245	1248	1249	rBV2	3076	2679	0.19%	0.016%
59	10.802	1266	1270	1278	rVB5	20661	35110	2.51%	0.205%
60	10.908	1280	1288	1295	rBV	289984	528122	37.73%	3.081%
61	11.031	1301	1309	1319	rBV	189771	369737	26.42%	2.157%
62	11.149	1324	1329	1336	rVB10	9869	21228	1.52%	0.124%
63	11.208	1336	1339	1341	rBV4	3121	3890	0.28%	0.023%
64	11.314	1352	1357	1363	rBV7	5427	11920	0.85%	0.070%
65	11.443	1374	1379	1381	rBV4	5484	9202	0.66%	0.054%
66	11.466	1381	1383	1385	rVV3	3246	3194	0.23%	0.019%
67	11.537	1391	1395	1397	rBV3	3885	5040	0.36%	0.029%
68	11.590	1402	1404	1411	rVB7	6540	10772	0.77%	0.063%
69	11.695	1419	1422	1424	rBV4	8397	10262	0.73%	0.060%
70	12.095	1486	1490	1495	rBV8	6882	14266	1.02%	0.083%
71	12.154	1495	1500	1504	rVB8	5350	10793	0.77%	0.063%

LSC Area Percent Report

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Misc :
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Instrument : BNA_G
ClientSampleId : CB7N5

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_G\METHODS\SOM-EPA-BG092920MA.M
Title : SVOA CALIBRATION

72	12.224	1505	1512	1517	rBV2	70410	128501	9.18%	0.750%
73	12.395	1536	1541	1546	rVB4	19215	34478	2.46%	0.201%
74	12.577	1569	1572	1573	rBV3	5703	6006	0.43%	0.035%
75	12.806	1609	1611	1612	rBV2	4730	2652	0.19%	0.015%
76	12.876	1618	1623	1627	rBV8	8405	17649	1.26%	0.103%
77	14.116	1829	1834	1839	rBV	450512	630817	45.07%	3.681%
78	14.163	1840	1842	1847	rVB	25751	27708	1.98%	0.162%
79	14.416	1879	1885	1895	rVB	456286	731716	52.28%	4.269%
80	14.516	1898	1902	1910	rVB2	36886	69102	4.94%	0.403%
81	14.721	1931	1937	1946	rVB2	488031	777691	55.56%	4.538%
82	14.903	1964	1968	1975	rVB2	51233	90606	6.47%	0.529%
83	15.450	2056	2061	2067	rBV	268761	397088	28.37%	2.317%
84	15.708	2099	2105	2113	rBV	591986	908625	64.92%	5.302%
85	15.814	2118	2123	2129	rBV	185636	275468	19.68%	1.607%
86	16.214	2186	2191	2200	rVB	368867	521366	37.25%	3.042%
87	16.384	2217	2220	2226	rVB2	67302	100373	7.17%	0.586%
88	16.725	2272	2278	2282	rBV	213238	311334	22.24%	1.817%
89	16.766	2282	2285	2289	rVB	52763	64287	4.59%	0.375%
90	16.983	2319	2322	2326	rBV3	28554	36142	2.58%	0.211%
91	17.465	2398	2404	2411	rVB	667950	992939	70.94%	5.794%
92	17.565	2415	2421	2428	rVB2	643407	1014682	72.49%	5.920%
93	18.346	2549	2554	2560	rBV	209125	311050	22.22%	1.815%
94	19.610	2765	2769	2778	rBV3	82353	143988	10.29%	0.840%
95	19.845	2803	2809	2813	rBV	971567	1399680	100.00%	8.167%
96	21.378	3066	3070	3081	rBV2	178225	329246	23.52%	1.921%
97	21.737	3125	3131	3138	rBV2	732076	1212329	86.61%	7.074%
98	24.774	3639	3648	3658	rVB	487326	1290220	92.18%	7.528%
99	25.003	3678	3687	3701	rVB2	444982	1337025	95.52%	7.801%
100	27.583	4118	4126	4135	rVB8	29665	83653	5.98%	0.488%

Sum of corrected areas: 17138634

LSC Report - Integrated Chromatogram

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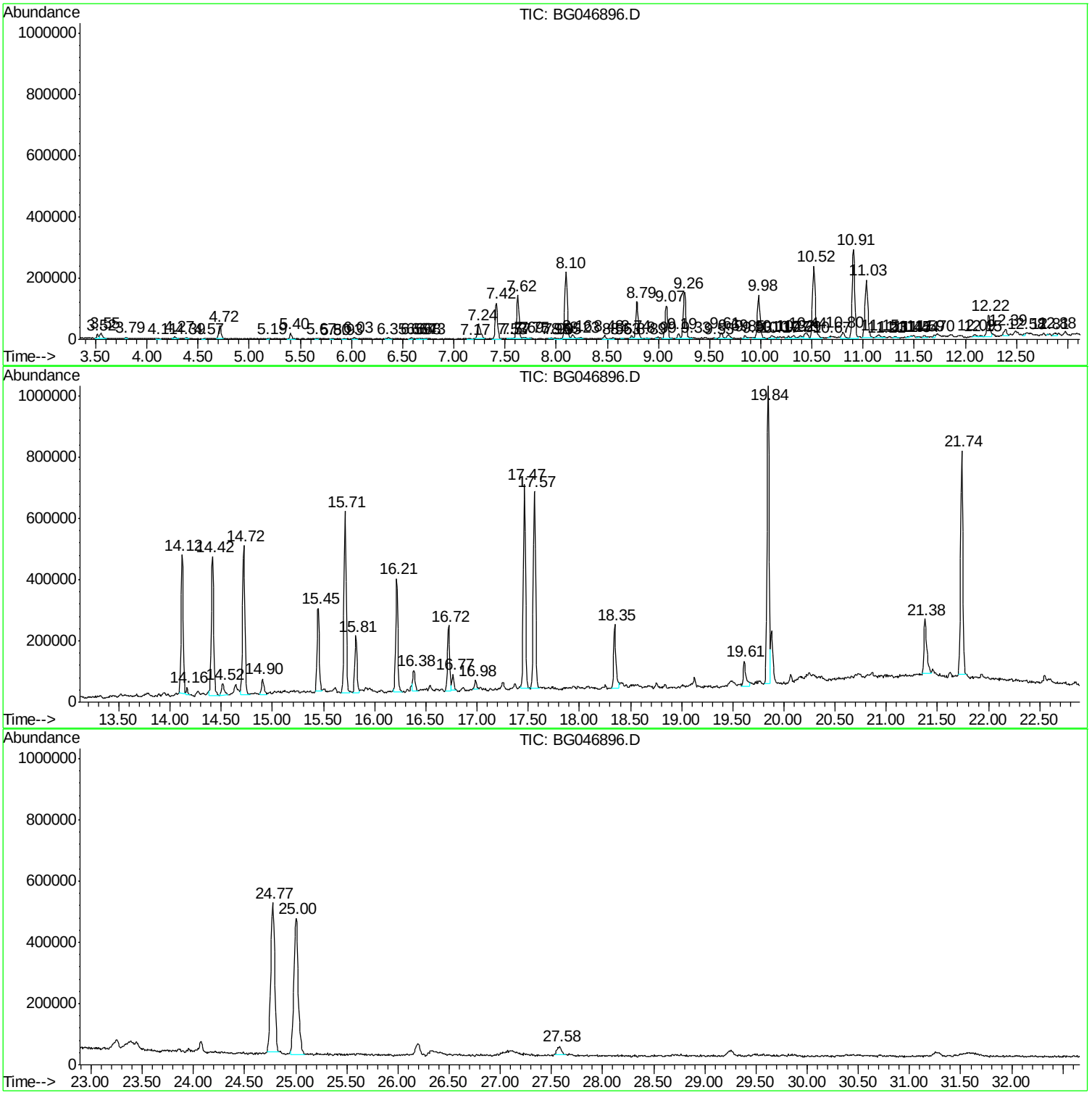
Data File : BG046896.D
Acq On : 14 Oct 2020 17:06
Operator : CG/JU
Sample : L4359-07
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : CB7N5

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG092920MA.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 G\DATA\BG101420\
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Instrument : BNA_G
ClientSampleId : CB7N5

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG092920MA.M
Quant Title : SVOA CALIBRATION

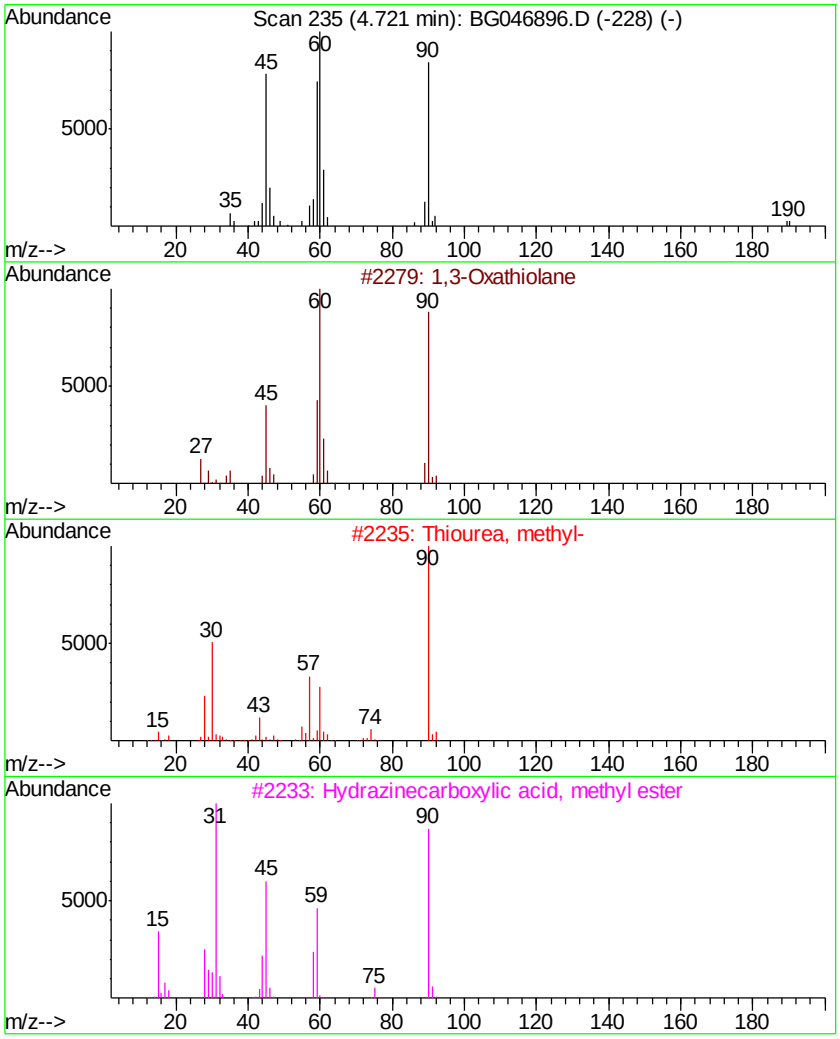
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

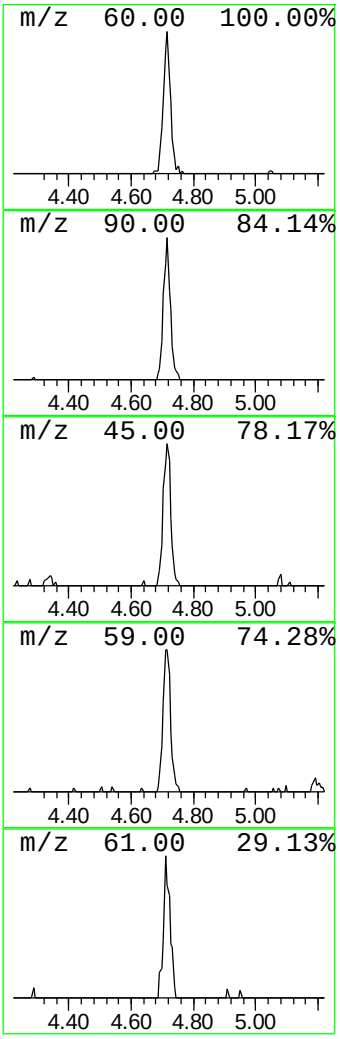
****

Peak Number 1 1,3-Oxathiolane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.72	3.73 ng/ul	69032	1,4-Dichlorobenzene-d4	8.10

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,3-Oxathiolane	90	C3H6OS	002094-97-5	64
2	Thiourea, methyl-	90	C2H6N2S	000598-52-7	50
3	Hydrazinecarboxylic acid, methyl...	90	C2H6N2O2	006294-89-9	45
4	1,3-Oxathiolane	90	C3H6OS	002094-97-5	40
5	3-Thietanol	90	C3H6OS	010304-16-2	40





Library Search Compound Report

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Sample : L4359-07
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : CB7N5

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG092920MA.M
Quant Title : SVOA CALIBRATION

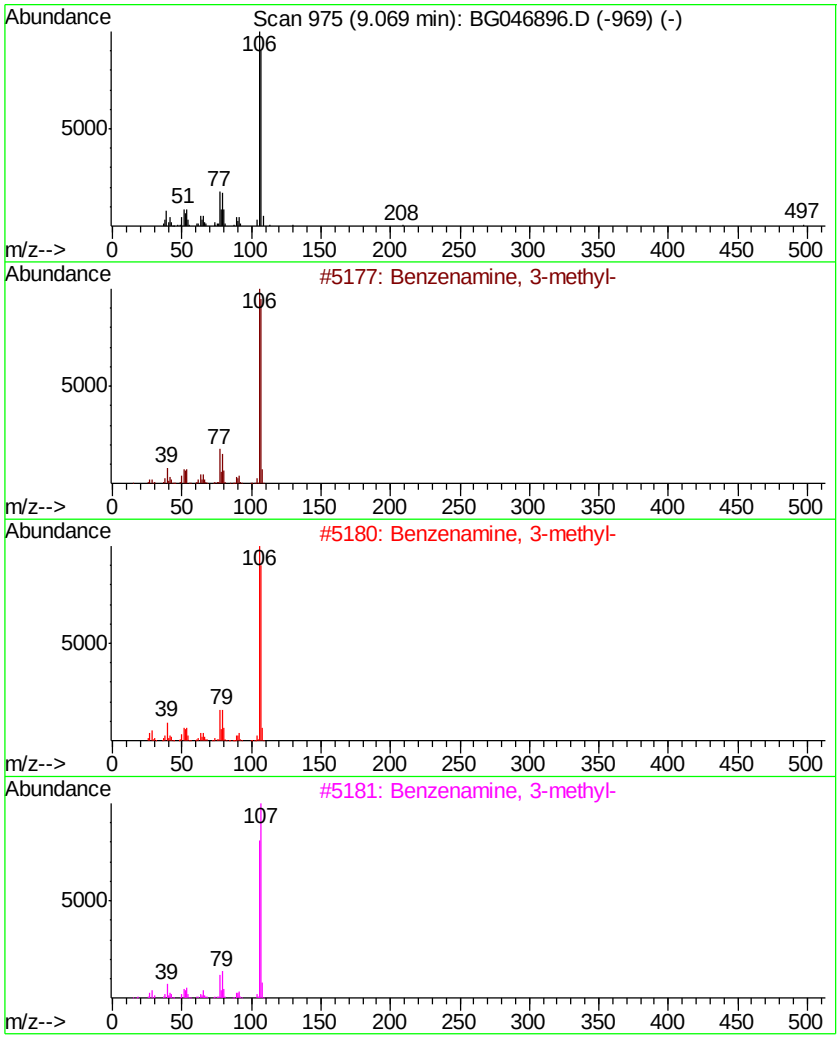
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

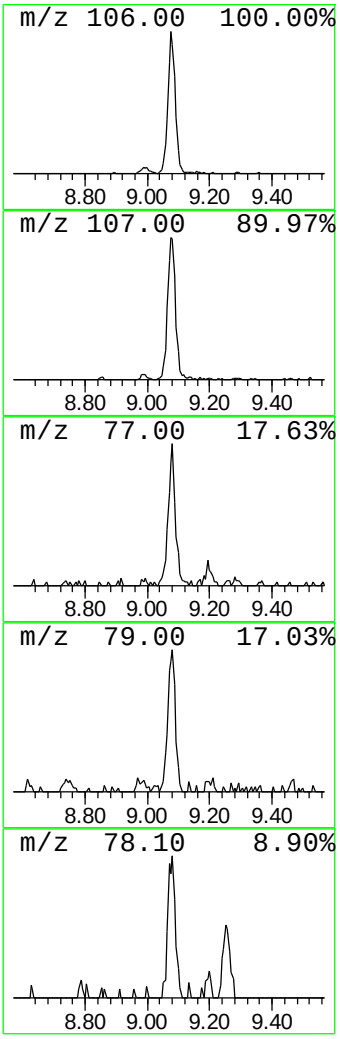
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.07	10.22 ng/ul	188890	1,4-Dichlorobenzene-d4	8.10

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzenamine, 3-methyl-	107	C7H9N	000108-44-1	97
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	94
5	Benzenamine, 3-methyl-	107	C7H9N	000108-44-1	94





Library Search Compound Report

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ClientSampleId : CB7N5

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG092920MA.M
Quant Title : SVOA CALIBRATION

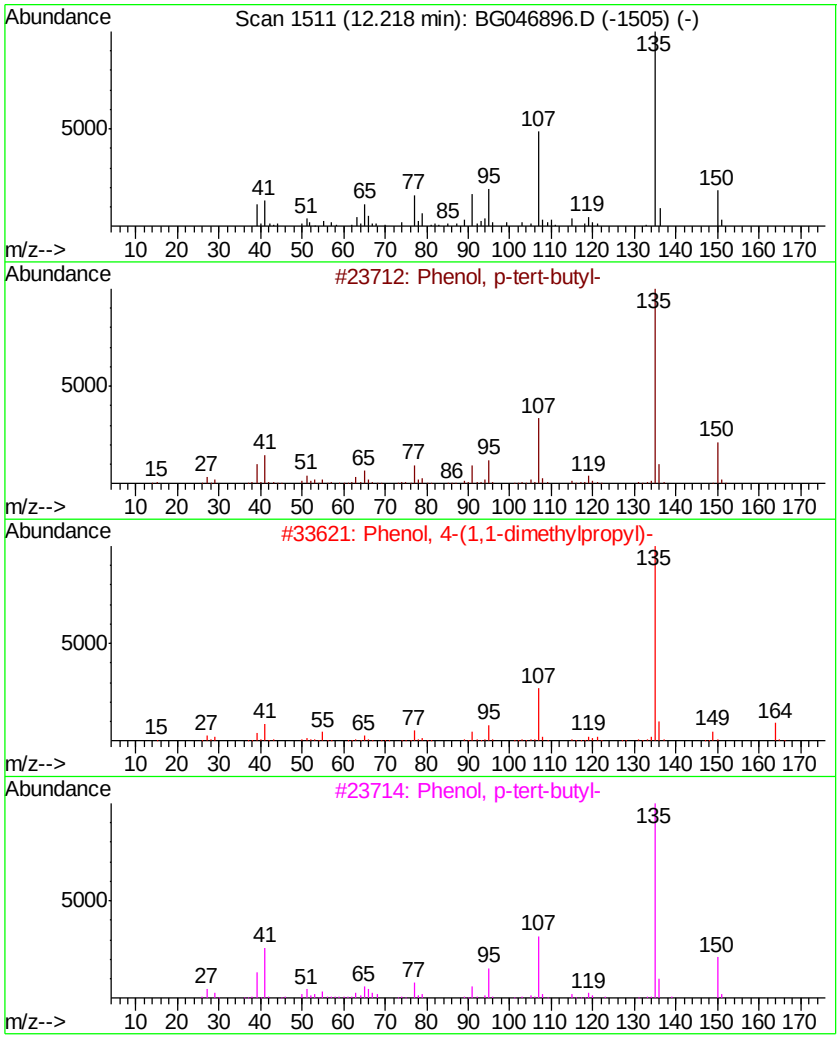
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

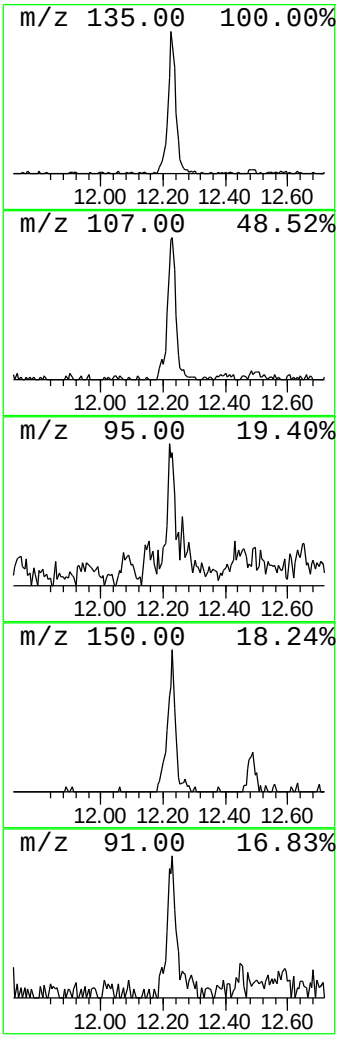
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.22	4.87 ng/ul	128501	Naphthalene-d8	10.91

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	91
2	Phenol, 4-(1,1-dimethylpropyl)-	164	C11H16O	000080-46-6	78
3	Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	72
4	Hexestrol	270	C18H22O2	005635-50-7	64
5	Phenol, m-tert-butyl-	150	C10H14O	000585-34-2	53





Library Search Compound Report

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ClientSampleId : CB7N5

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG092920MA.M
Quant Title : SVOA CALIBRATION

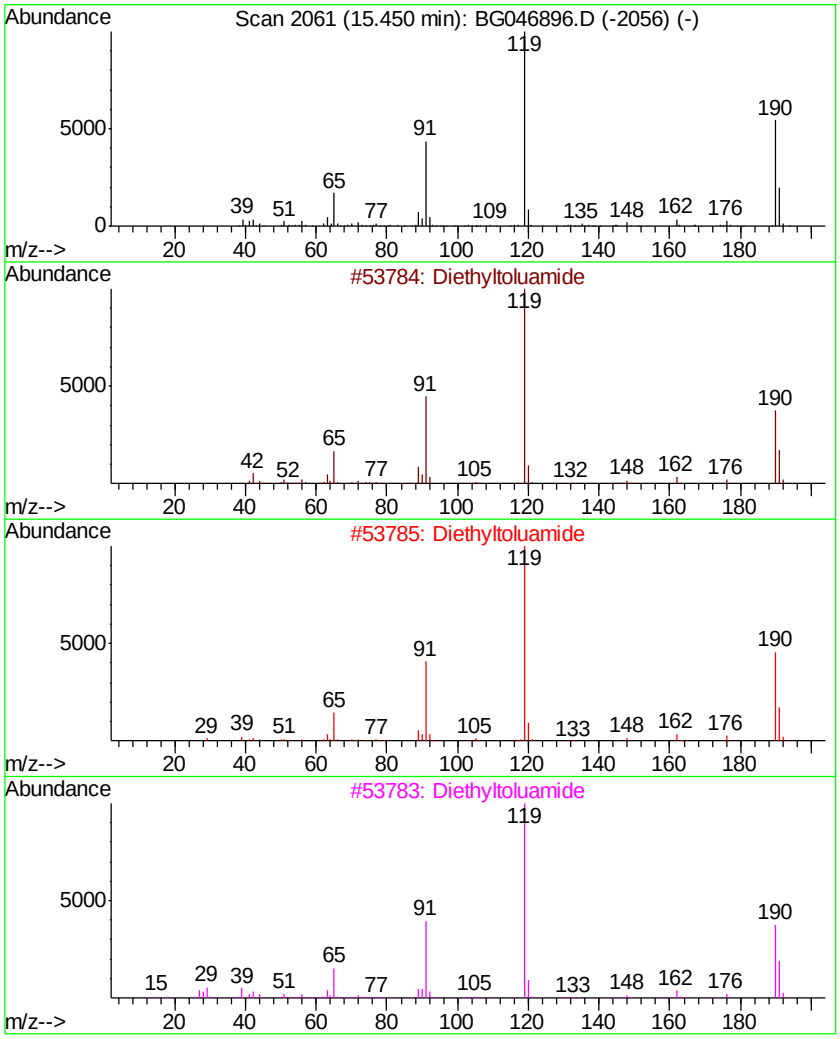
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

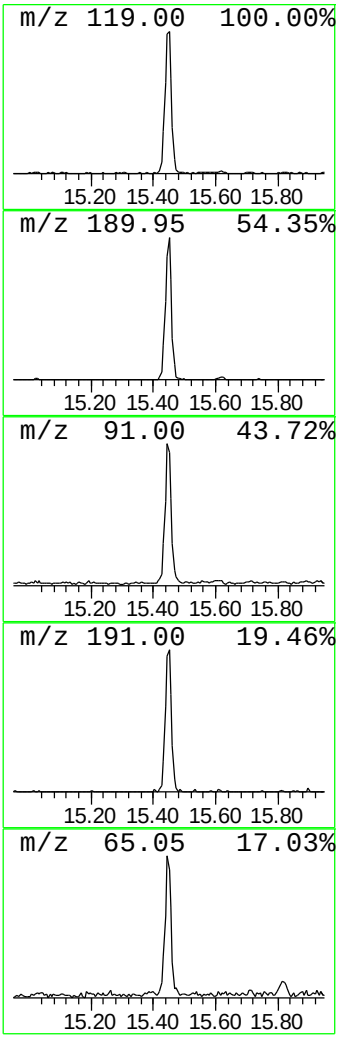
****

Peak Number 4 Diethyltoluamide Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.45	10.21 ng/ul	397088	Acenaphthene-d10	14.72

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Diethyltoluamide	191	C12H17NO	000134-62-3	94
2	Diethyltoluamide	191	C12H17NO	000134-62-3	94
3	Diethyltoluamide	191	C12H17NO	000134-62-3	91
4	Benzamide, N,N-diethyl-4-methyl-	191	C12H17NO	002728-05-4	87
5	Diethyltoluamide	191	C12H17NO	000134-62-3	72





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Sample : L4359-07
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : CB7N5

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG092920MA.M
Quant Title : SVOA CALIBRATION

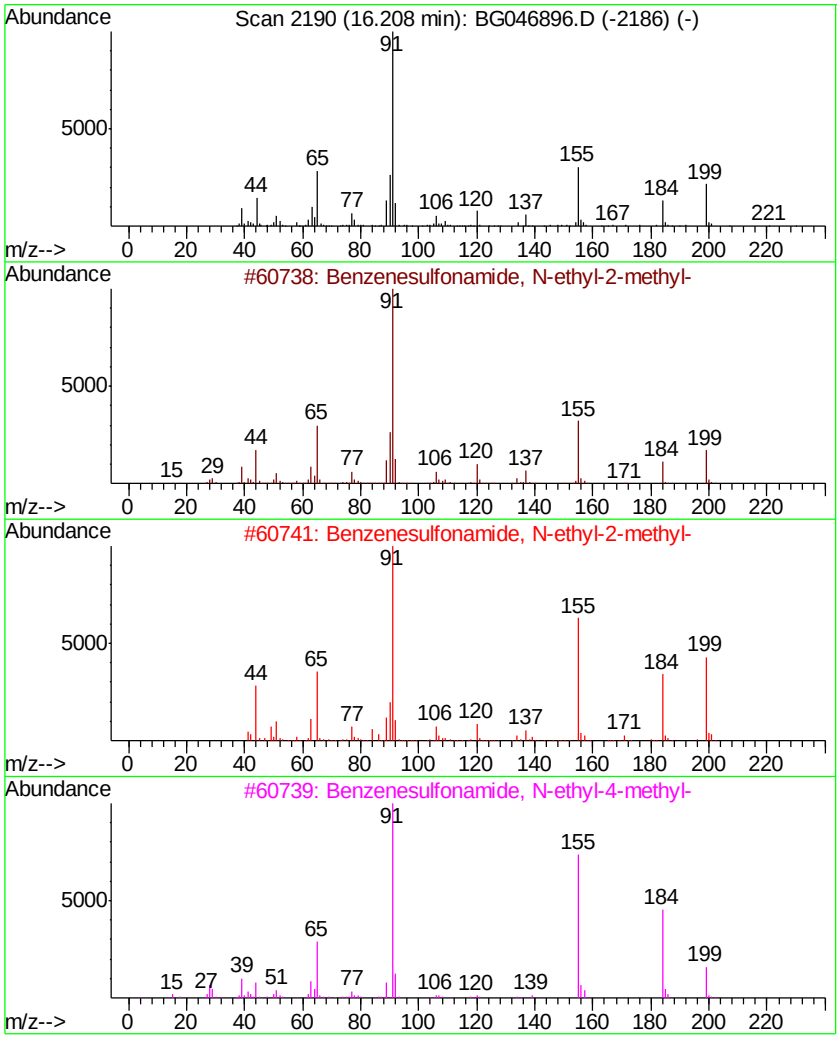
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

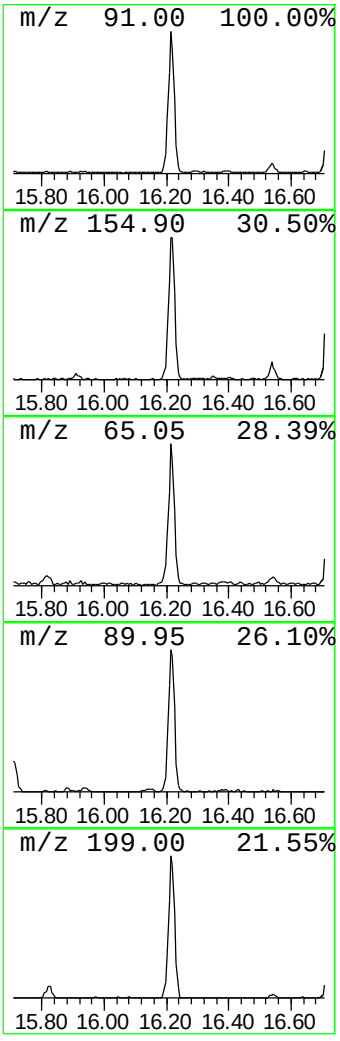
****

Peak Number 5 Benzenesulfonamide, N-ethyl... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.21	10.50 ng/ul	521366	Phenanthrene-d10	17.47

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzenesulfonamide, N-ethyl-2-me...	199	C9H13NO2S	001077-56-1	91
2	Benzenesulfonamide, N-ethyl-2-me...	199	C9H13NO2S	001077-56-1	58
3	Benzenesulfonamide, N-ethyl-4-me...	199	C9H13NO2S	000080-39-7	49
4	p-Toluenesulfonylacetonitrile	195	C9H9NO2S	005697-44-9	47
5	4-Toluenesulfonylmethyl isocyanide	195	C9H9NO2S	036635-61-7	47





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Data File : BG046896.D
Acq On : 14 Oct 2020 17:06
Operator : CG/JU
Sample : L4359-07
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : CB7N5

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG092920MA.M
Quant Title : SVOA CALIBRATION

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

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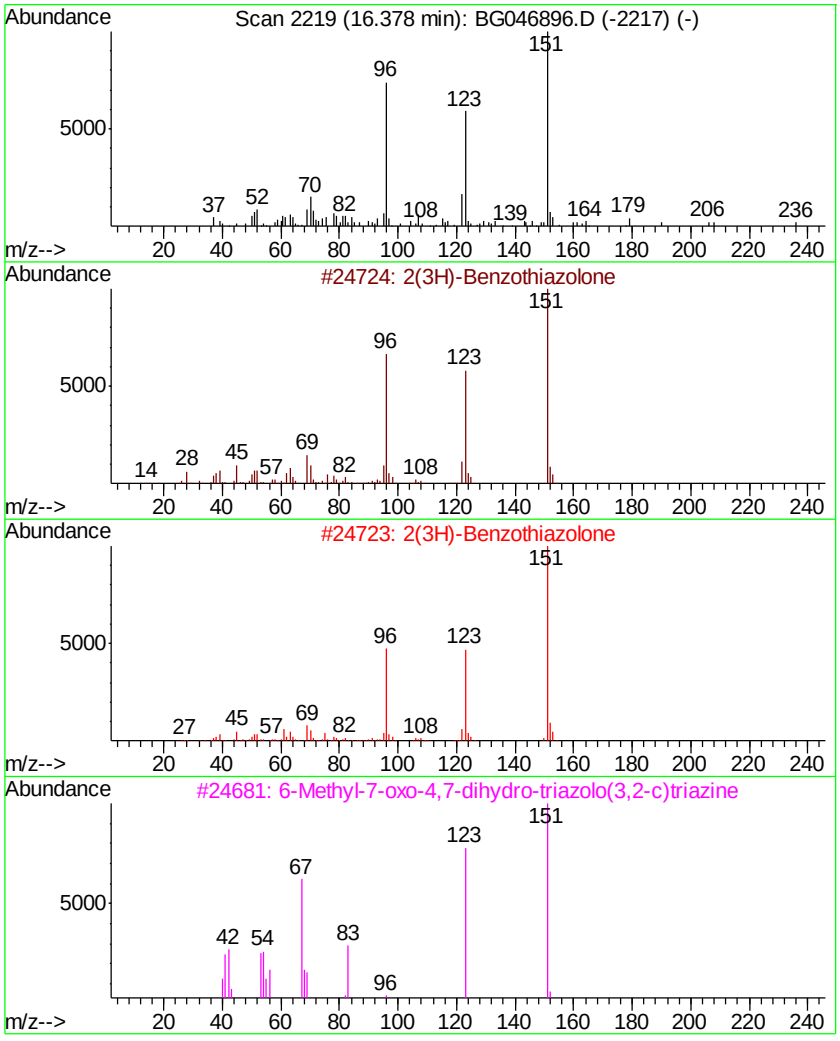
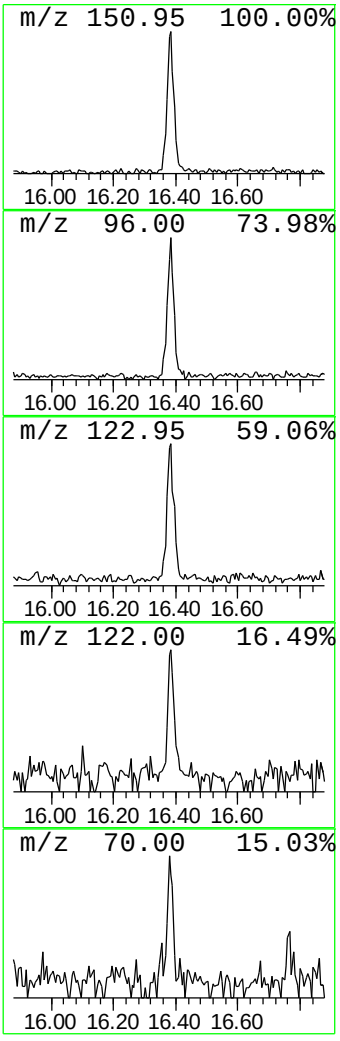
Peak Number 6 2(3H)-Benzothiazolone Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.38	2.02 ng/ul	100373	Phenanthrene-d10	17.47

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2(3H)-Benzothiazolone	151	C7H5NOS	000934-34-9	86
2	2(3H)-Benzothiazolone	151	C7H5NOS	000934-34-9	58
3	6-Methyl-7-oxo-4,7-dihydro-triaz...	151	C5H5N5O	057250-39-2	38
4	Thieno[2,3-b]pyridine-N-oxide	151	C7H5NOS	025557-50-0	32
5	2(3H)-Benzoxazolethione	151	C7H5NOS	002382-96-9	27





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Data File : BG046896.D
Acq On : 14 Oct 2020 17:06
Operator : CG/JU
Sample : L4359-07
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : CB7N5

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG092920MA.M
Quant Title : SVOA CALIBRATION

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

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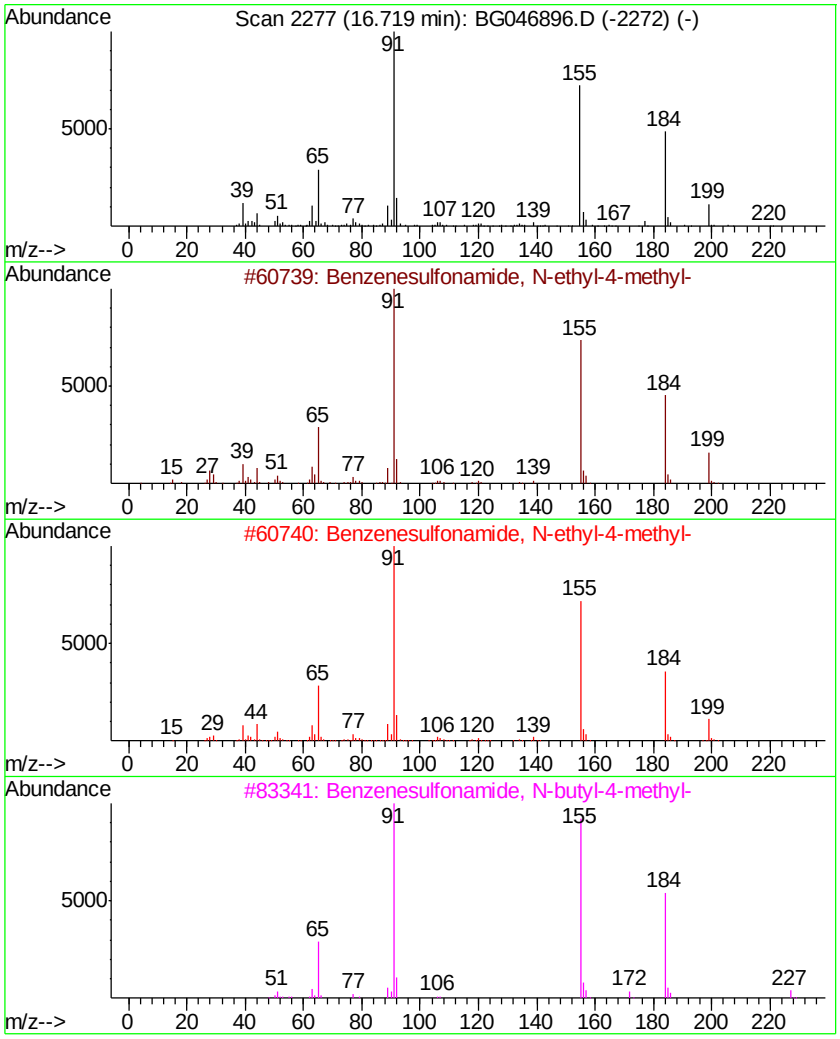
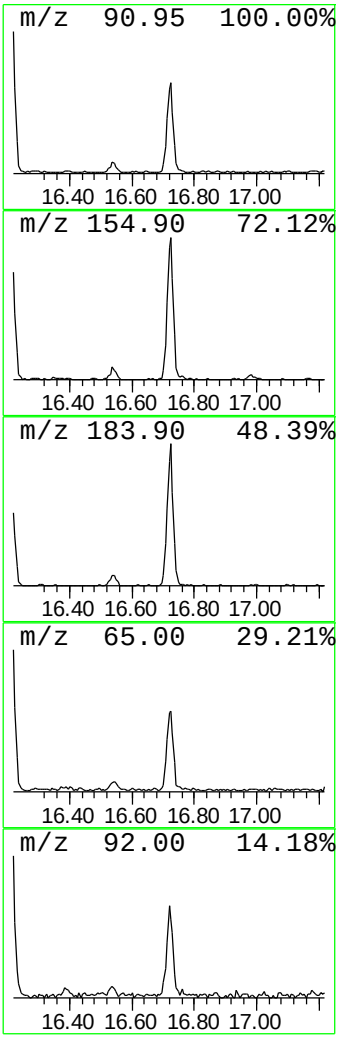
Peak Number 7 Benzenesulfonamide, N-ethyl... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.72	6.27 ng/ul	311334	Phenanthrene-d10	17.47

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	Benzenesulfonamide, N-butyl-4-me...	227	C11H17NO2S	001907-65-9	86
4	N-Phenethyl-p-toluenesulfonamide	275	C15H17NO2S	005450-75-9	78
5	Furazan-3-carboxylic acid, 4-ami...	325	C12H15N5O4S	1000304-69-4	78





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Acq On : 14 Oct 2020 17:06
Operator : CG/JU
Sample : L4359-07
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : CB7N5

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG092920MA.M
Quant Title : SVOA CALIBRATION

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

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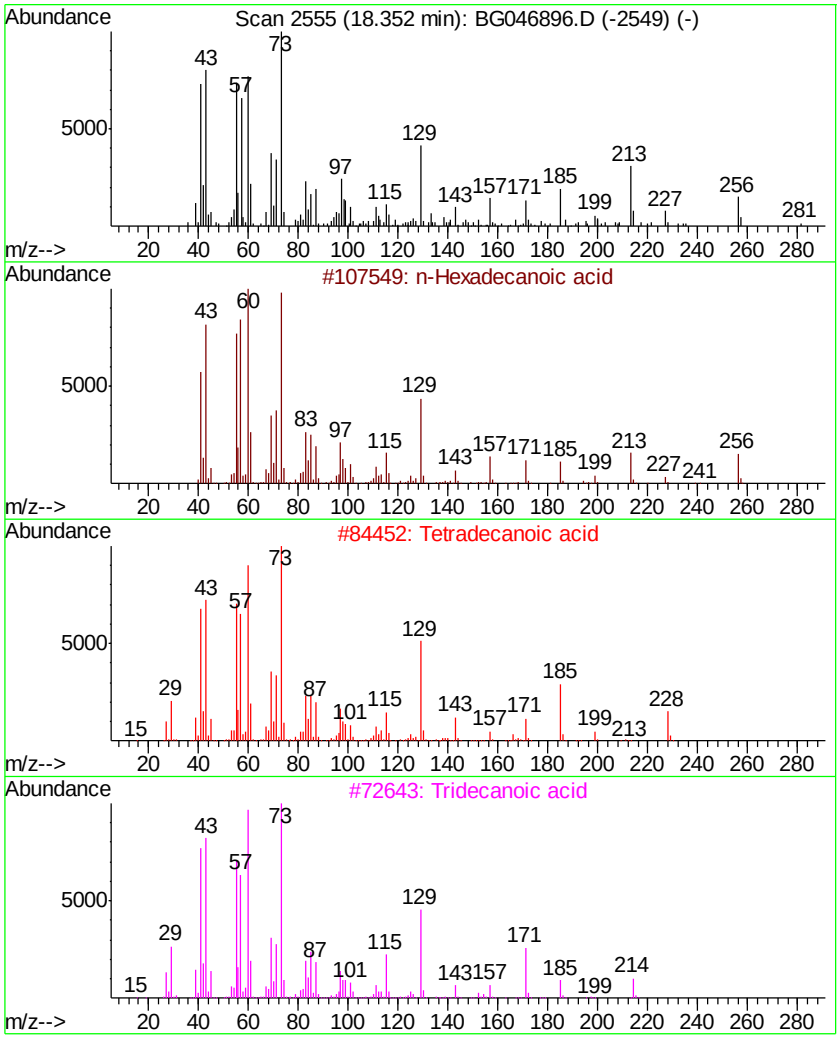
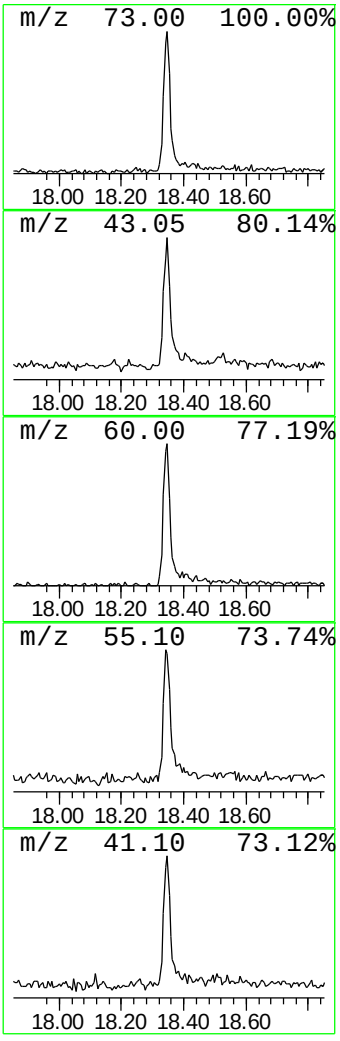
Peak Number 8 n-Hexadecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.35	6.27 ng/ul	311050	Phenanthrene-d10	17.47

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2	Tetradecanoic acid	228	C14H28O2	000544-63-8	91
3	Tridecanoic acid	214	C13H26O2	000638-53-9	90
4	Pentadecanoic acid	242	C15H30O2	001002-84-2	87
5	Tetradecanoic acid	228	C14H28O2	000544-63-8	80





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Data File : BG046896.D
Acq On : 14 Oct 2020 17:06
Operator : CG/JU
Sample : L4359-07
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : CB7N5

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG092920MA.M
Quant Title : SVOA CALIBRATION

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

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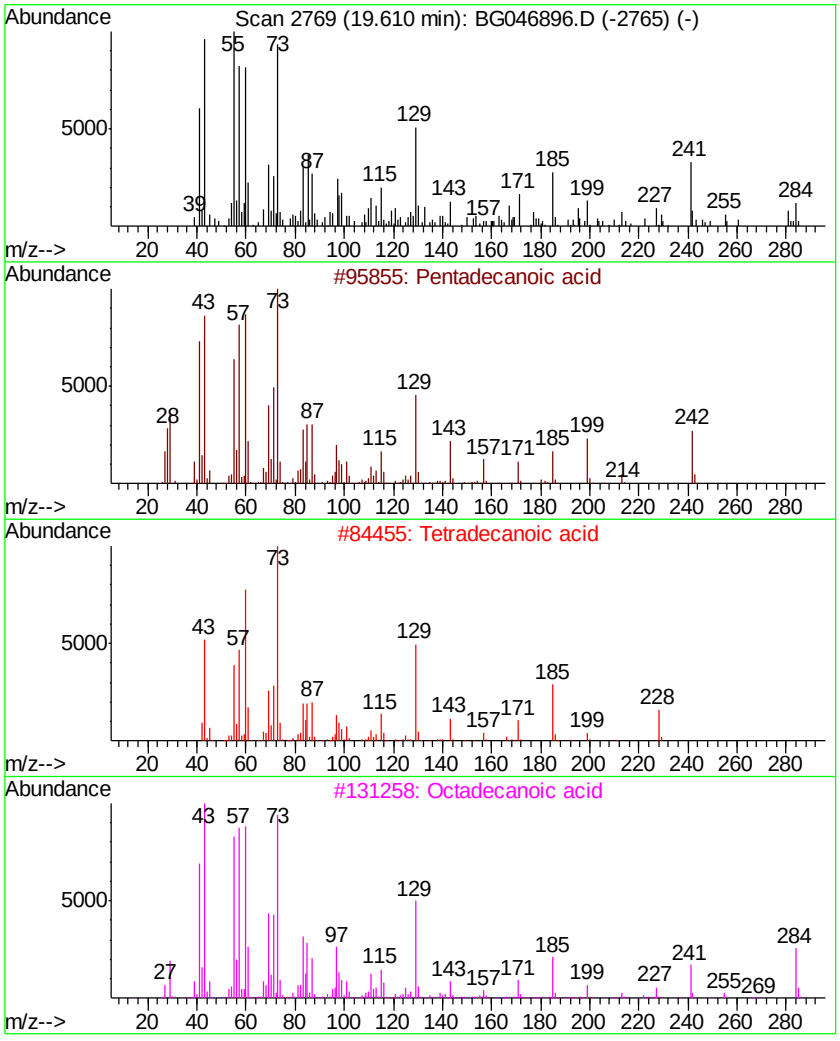
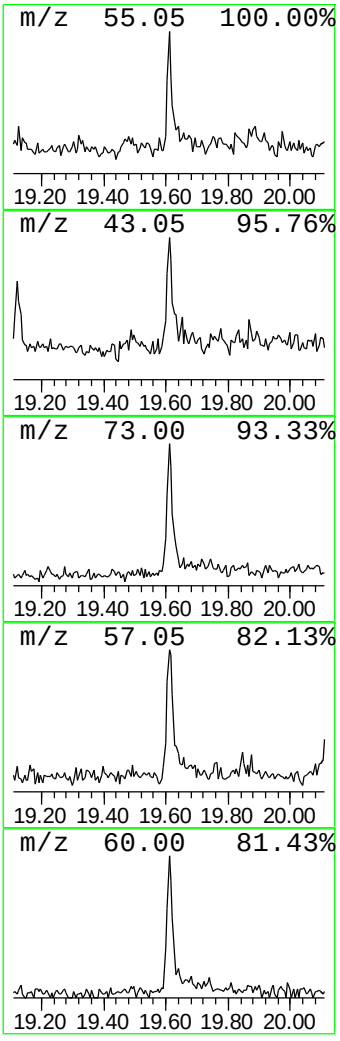
Peak Number 9 Pentadecanoic acid Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.61	2.38 ng/ul	143988	Chrysene-d12	21.74

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentadecanoic acid	242	C15H30O2	001002-84-2	80
2	Tetradecanoic acid	228	C14H28O2	000544-63-8	72
3	Octadecanoic acid	284	C18H36O2	000057-11-4	72
4	Tetradecanoic acid	228	C14H28O2	000544-63-8	64
5	Tridecanoic acid	214	C13H26O2	000638-53-9	59





Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG101420\
Data File : BG046896.D
Acq On : 14 Oct 2020 17:06
Operator : CG/JU
Sample : L4359-07
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : CB7N5

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG092920MA.M
Quant Title : SVOA CALIBRATION

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

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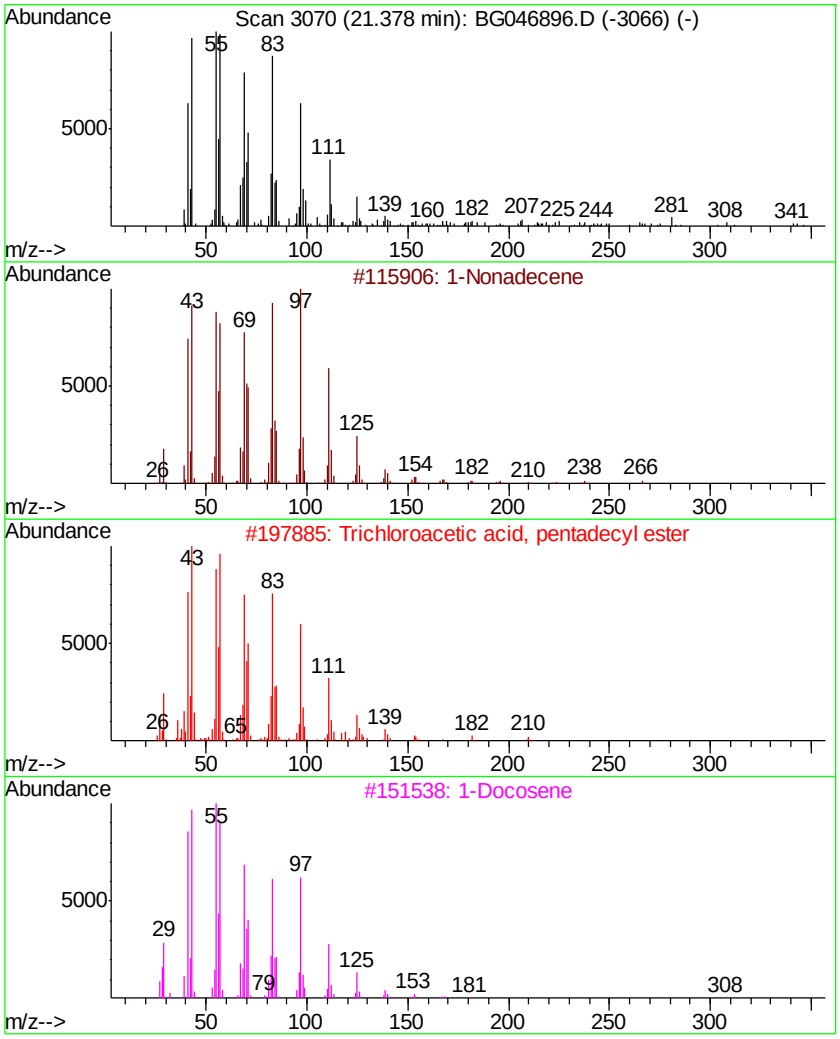
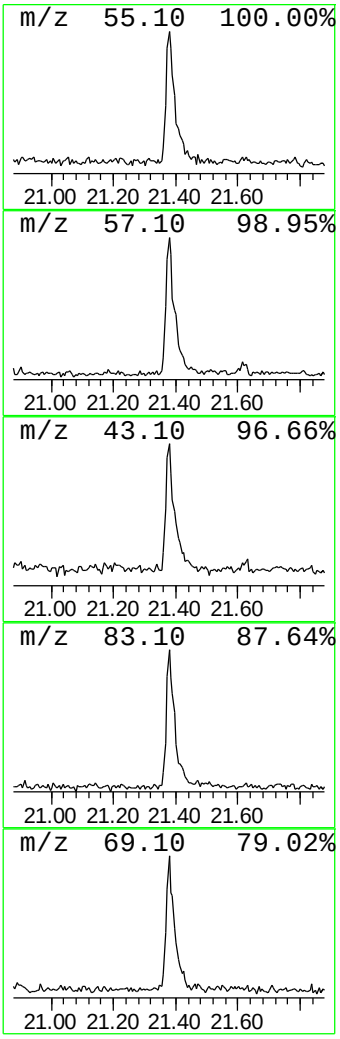
Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.38	5.43 ng/ul	329246	Chrysene-d12	21.74

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Nonadecene	266	C19H38	018435-45-5	94
2	Trichloroacetic acid, pentadecyl...	372	C17H31Cl3O2	074339-53-0	94
3	1-Docosene	308	C22H44	001599-67-3	93
4	Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	91
5	Bromoacetic acid, hexadecyl ester	362	C18H35BrO2	005454-48-8	90





Tentatively Identified Compound (LSC) summary

Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG101420\
Data File : BG046896.D
Acq On : 14 Oct 2020 17:06
Operator : CG/JU
Sample : L4359-07
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : CB7N5

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG092920MA.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	4.72	3.7	ng/ul	69032	1	8.10	369796	20.0
Benzenamine, 3-me...	9.07	10.2	ng/ul	188890	1	8.10	369796	20.0
Phenol, p-tert-bu...	12.22	4.9	ng/ul	128501	2	10.91	528122	20.0
Diethyltoluamide	15.45	10.2	ng/ul	397088	3	14.72	777691	20.0
Benzenesulfonamid...	16.21	10.5	ng/ul	521366	4	17.47	992939	20.0
2(3H)-Benzothiazo...	16.38	2.0	ng/ul	100373	4	17.47	992939	20.0
Benzenesulfonamid...	16.72	6.3	ng/ul	311334	4	17.47	992939	20.0
n-Hexadecanoic acid	18.35	6.3	ng/ul	311050	4	17.47	992939	20.0
Pentadecanoic acid	19.61	2.4	ng/ul	143988	5	21.74	1212330	20.0
(DEL) Alkane: Str...	21.38	5.4	ng/ul	329246	5	21.74	1212330	20.0