LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111723\
Data File : BP018251.D
Acq On : 18 Nov 2023 09:17
Operator : MA/JU
Sample : 05370-03
Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : GCDP9

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_P\Methods\SFAM-EPA-BP111023.MA.M
Title : SVOA CALIBRATION

Signal : TIC: BP018251.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.158	9	12	16	rBV	15842	19106	0.75%	0.081%
2	3.770	113	116	120	rBV5	6336	9158	0.36%	0.039%
3	3.899	134	138	144	rBV6	3922	9025	0.35%	0.038%
4	4.246	195	197	203	rVB6	2984	5361	0.21%	0.023%
5	5.658	432	437	440	rBV6	1931	4031	0.16%	0.017%
6	6.728	615	619	625	rVB3	5923	11476	0.45%	0.049%
7	6.958	650	658	672	rBV	87088	196765	7.73%	0.838%
8	7.128	681	687	700	rBV	233798	385886	15.16%	1.644%
9	7.299	710	716	738	rBV	312959	561618	22.07%	2.392%
10	7.775	790	797	815	rBV	263608	480131	18.87%	2.045%
11	7.887	815	816	821	rVB4	2173	2772	0.11%	0.012%
12	8.505	915	921	943	rBV	239891	496105	19.50%	2.113%
13	8.675	949	950	956	rVB6	2222	2649	0.10%	0.011%
14	8.987	996	1003	1016	rBV	294957	546048	21.46%	2.326%
15	9.152	1025	1031	1036	rVB	32285	51804	2.04%	0.221%
16	9.387	1069	1071	1078	rVB7	1625	2656	0.10%	0.011%
17	9.699	1117	1124	1139	rBV	271763	521767	20.50%	2.222%
18	10.228	1208	1214	1241	rBV	260569	702558	27.61%	2.993%
19	10.493	1257	1259	1262	rVB4	2215	2687	0.11%	0.011%
20	10.593	1268	1276	1287	rBV	375603	617177	24.25%	2.629%
21	10.763	1300	1305	1310	rBV9	2695	5182	0.20%	0.022%
22	10.834	1310	1317	1322	rBV8	13261	36761	1.44%	0.157%
23	10.981	1334	1342	1354	rVB8	17083	47351	1.86%	0.202%
24	11.087	1354	1360	1361	rBV6	3142	4685	0.18%	0.020%
25	11.122	1363	1366	1367	rVV3	7454	7345	0.29%	0.031%
26	11.146	1367	1370	1380	rVB9	9185	20032	0.79%	0.085%
27	11.322	1394	1400	1404	rBV8	3538	6695	0.26%	0.029%
28	11.405	1404	1414	1423	rBV4	27929	85509	3.36%	0.364%
29	11.481	1423	1427	1431	rBV3	8205	13950	0.55%	0.059%
30	11.528	1431	1435	1439	rVV3	10394	16095	0.63%	0.069%
31	11.646	1450	1455	1461	rVB2	25946	42474	1.67%	0.181%
32	12.240	1550	1556	1562	rBV8	3927	7985	0.31%	0.034%
33	12.646	1617	1625	1626	rBV8	1777	4100	0.16%	0.017%
34	12.757	1643	1644	1652	rVV8	2181	2837	0.11%	0.012%
35	13.022	1686	1689	1694	rVB6	4259	5816	0.23%	0.025%
36	13.299	1727	1736	1744	rBV2	48824	82705	3.25%	0.352%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111723\
Data File : BP018251.D
Acq On : 18 Nov 2023 09:17
Operator : MA/JU
Sample : 05370-03
Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCDP9

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_P\Methods\SFAM-EPA-BP111023.MA.M
Title : SVOA CALIBRATION

37	13.669	1791	1799	1804	rBV6	5376	11573	0.45%	0.049%
38	13.734	1804	1810	1818	rBV2	9165	23745	0.93%	0.101%
39	13.881	1829	1835	1850	rBV	810855	1179642	46.36%	5.025%
40	14.157	1876	1882	1899	rBV	845070	1233186	48.46%	5.253%
41	14.457	1926	1933	1944	rBV	551664	805598	31.66%	3.431%
42	14.775	1982	1987	1995	rBV4	27867	87383	3.43%	0.372%
43	15.104	2038	2043	2047	rBV4	5463	8604	0.34%	0.037%
44	15.204	2055	2060	2063	rBV	13368	17877	0.70%	0.076%
45	15.234	2063	2065	2072	rVB8	2799	5059	0.20%	0.022%
46	15.322	2077	2080	2084	rVB6	2807	4445	0.17%	0.019%
47	15.463	2098	2104	2112	rBV	1147645	1606916	63.15%	6.845%
48	15.534	2112	2116	2123	rVB6	14390	26377	1.04%	0.112%
49	15.628	2126	2132	2144	rBV	318323	567544	22.30%	2.417%
50	15.804	2160	2162	2165	rVB4	2531	2936	0.12%	0.013%
51	15.881	2170	2175	2178	rBV	19899	26550	1.04%	0.113%
52	15.922	2178	2182	2185	rVV4	10580	15029	0.59%	0.064%
53	15.963	2186	2189	2192	rVB4	4694	6261	0.25%	0.027%
54	16.122	2211	2216	2221	rVB9	3123	5545	0.22%	0.024%
55	16.181	2221	2226	2237	rBV	59374	83330	3.27%	0.355%
56	16.275	2237	2242	2246	rVV	159551	212727	8.36%	0.906%
57	16.334	2246	2252	2258	rVV3	194192	393942	15.48%	1.678%
58	16.398	2258	2263	2267	rVV3	153793	277851	10.92%	1.183%
59	16.440	2267	2270	2275	rVV	104450	138490	5.44%	0.590%
60	16.498	2275	2280	2281	rVV2	55454	80455	3.16%	0.343%
61	16.540	2285	2287	2292	rVV	54183	74278	2.92%	0.316%
62	16.598	2292	2297	2304	rVV3	121831	238654	9.38%	1.017%
63	16.669	2304	2309	2313	rVV	150492	222651	8.75%	0.948%
64	16.710	2313	2316	2318	rVV3	46942	66818	2.63%	0.285%
65	16.740	2318	2321	2330	rVB	74117	113334	4.45%	0.483%
66	16.881	2341	2345	2348	rBV3	7086	9605	0.38%	0.041%
67	16.910	2348	2350	2355	rVV6	4500	5737	0.23%	0.024%
68	16.957	2355	2358	2362	rVV6	6129	9449	0.37%	0.040%
69	16.998	2362	2365	2369	rVB3	8049	11839	0.47%	0.050%
70	17.045	2369	2373	2375	rBV4	4307	6482	0.25%	0.028%
71	17.081	2376	2379	2382	rVB5	5175	5656	0.22%	0.024%
72	17.116	2383	2385	2388	rVB4	4741	3466	0.14%	0.015%
73	17.157	2390	2392	2394	rBV3	2582	3061	0.12%	0.013%
74	17.240	2400	2406	2416	rBV	724523	1050638	41.29%	4.475%
75	17.345	2417	2424	2440	rVV	1264290	1864318	73.26%	7.941%
76	17.504	2447	2451	2457	rVB4	9833	17253	0.68%	0.073%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111723\
Data File : BP018251.D
Acq On : 18 Nov 2023 09:17
Operator : MA/JU
Sample : 05370-03
Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCDP9

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_P\Methods\SFAM-EPA-BP111023.MA.M
Title : SVOA CALIBRATION

77	17.875	2510	2514	2519	rBV8	14232	16488	0.65%	0.070%
78	18.098	2547	2552	2560	rBV2	56608	97929	3.85%	0.417%
79	18.187	2564	2567	2573	rVB3	12011	17504	0.69%	0.075%
80	18.310	2583	2588	2593	rBV	95766	140645	5.53%	0.599%
81	19.175	2732	2735	2741	rVB5	15499	22538	0.89%	0.096%
82	19.339	2759	2763	2769	rBV2	66700	93877	3.69%	0.400%
83	19.604	2802	2808	2816	rBV	1760029	2387850	93.84%	10.171%
84	21.380	3105	3110	3121	rBV	913559	1266716	49.78%	5.396%
85	23.692	3495	3503	3515	rVB2	1254817	2544703	100.00%	10.839%
86	23.851	3524	3530	3545	rVB	611140	1346353	52.91%	5.735%

Sum of corrected areas: 23477209

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111723\
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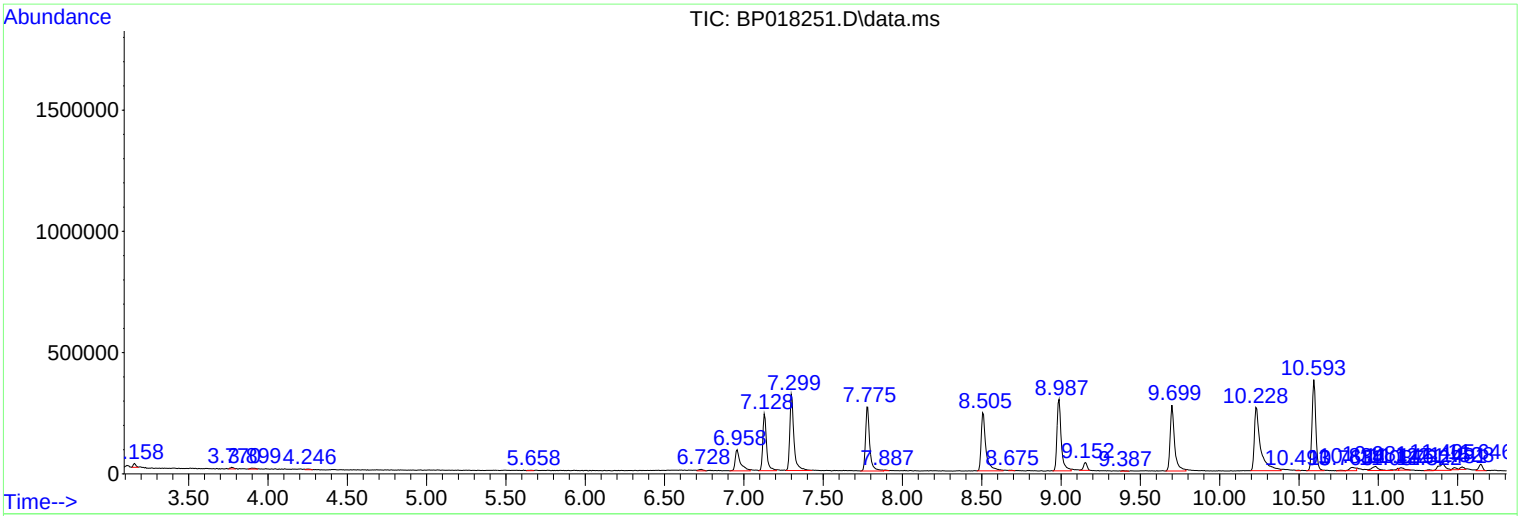
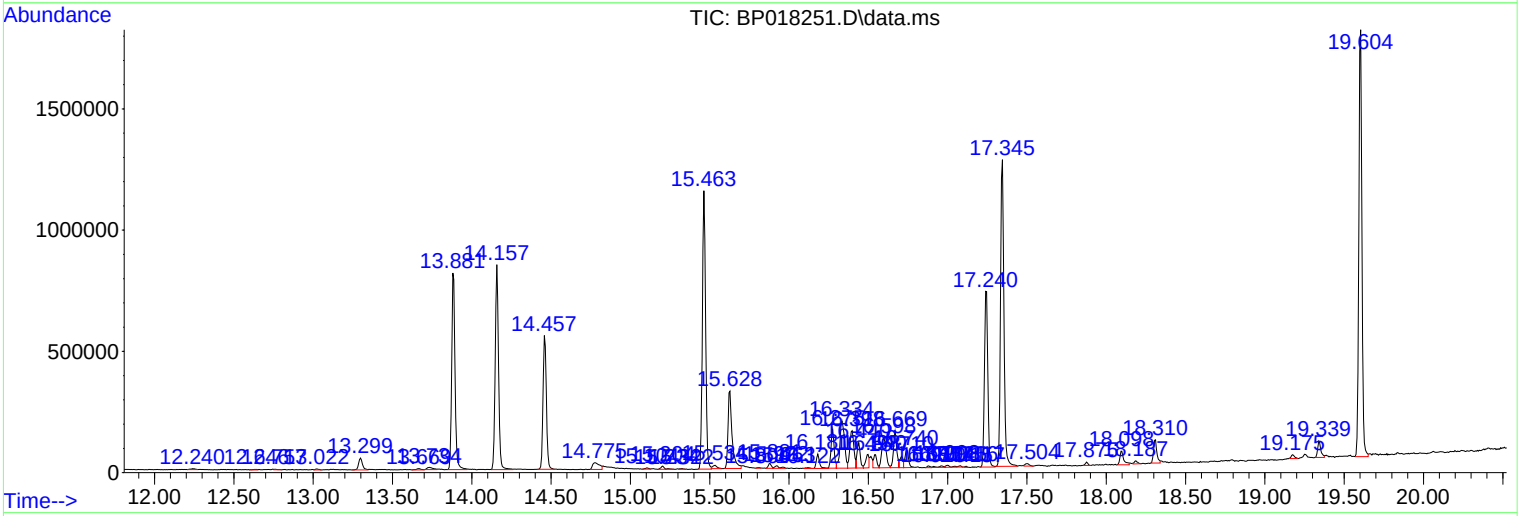
Operator : MA/JU
Sample : 05370-03
Misc :
ALS Vial : 40 Sample Multiplier: 1

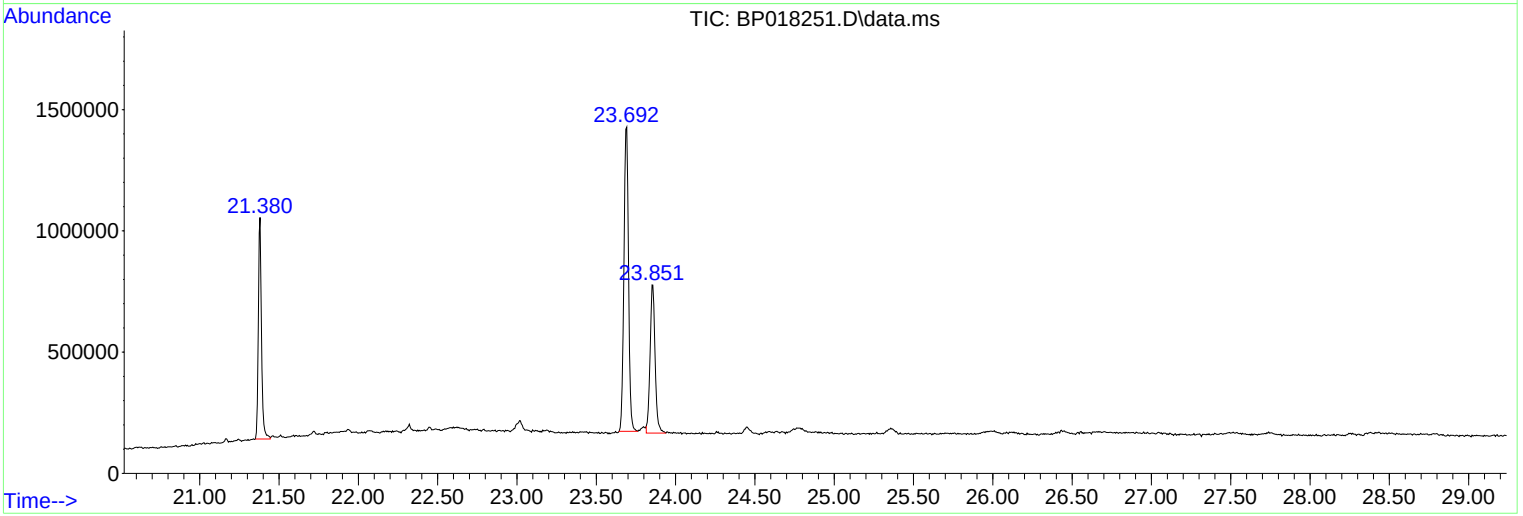
Instrument :
BNA_P
ClientSampleId :
GCDP9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P







Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111723\
Data File : BP018251.D
Acq On : 18 Nov 2023 09:17
Operator : MA/JU
Sample : 05370-03
Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCDP9

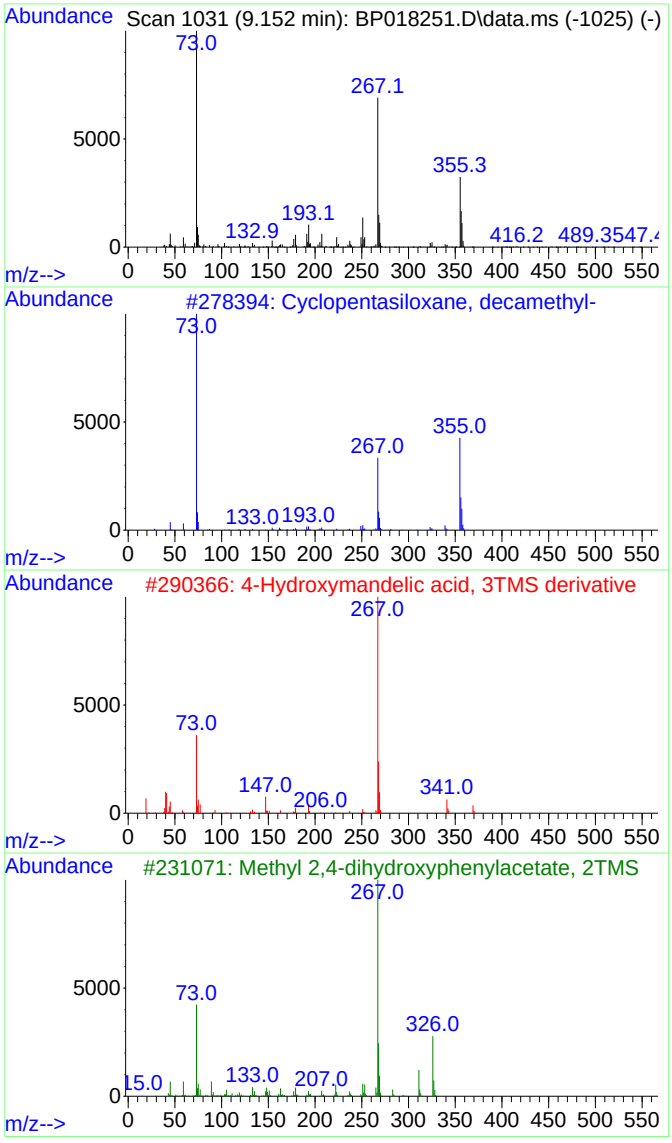
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
Quant Title : SVOA CALIBRATION

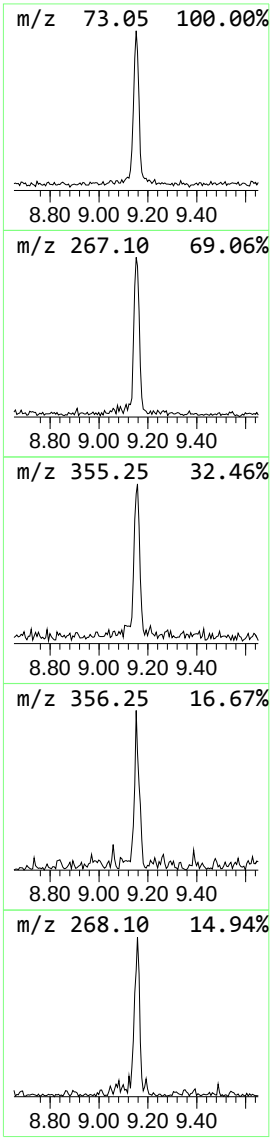
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 Cyclopentasiloxane, decamet... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.152	2.16 ng/ul	51804	1,4-Dichlorobenzene-d4	7.781

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentasiloxane, decamethyl-	370	C10H30O5Si5	000541-02-6	78
2	4-Hydroxymandelic acid, 3TMS der...	384	C17H32O4Si3	037148-64-4	47
3	Methyl 2,4-dihydroxyphenylacetat...	326	C15H26O4Si2	1010497-69-7	47
4	[(4-Hexylbenzene-1,3-diyl)bis(ox...	338	C18H34O2Si2	134273-01-1	47
5	2-Amino-4-methylphenol, 2TMS der...	267	C13H25NOSi2	1000373-28-1	47





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111723\
Data File : BP018251.D
Acq On : 18 Nov 2023 09:17
Operator : MA/JU
Sample : 05370-03
Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCDP9

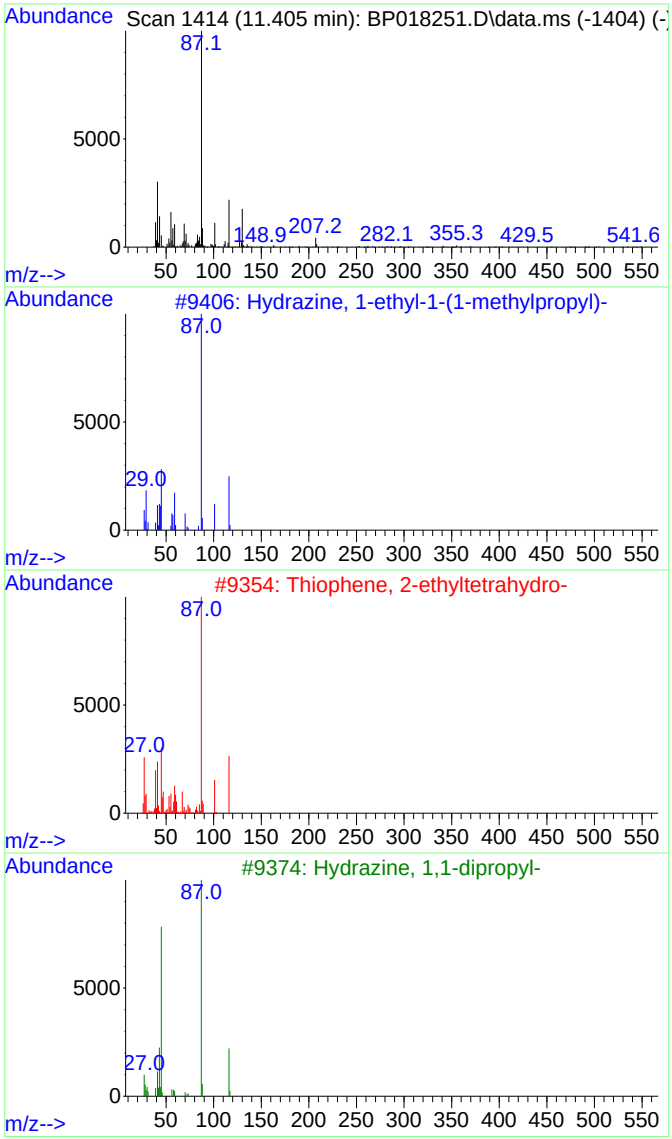
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
Quant Title : SVOA CALIBRATION

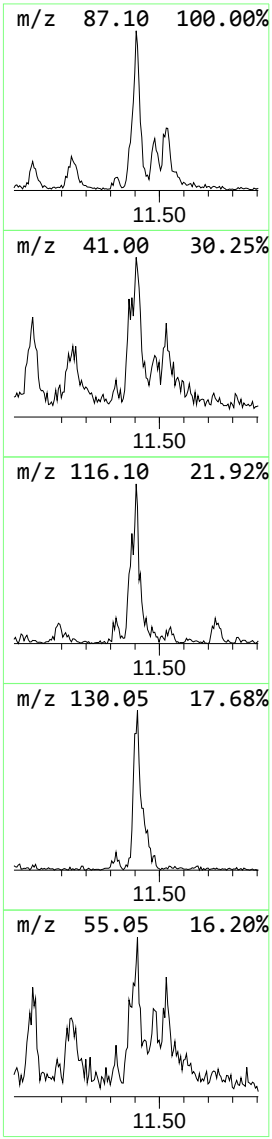
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 Hydrazine, 1-ethyl-1-(1-met... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.405	2.77 ng/ul	85509	Naphthalene-d8	10.593

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hydrazine, 1-ethyl-1-(1-methylpr...	116	C6H16N2	020325-97-7	64
2	Thiophene, 2-ethyltetrahydro-	116	C6H12S	001551-32-2	64
3	Hydrazine, 1,1-dipropyl-	116	C6H16N2	004986-50-9	52
4	2-[2-[2-[2-(2-Butoxyethoxy)ethox...	336	C16H32O7	1000351-94-0	50
5	4-(4-Chloro-2-methylphenoxy)buty...	438	C26H43ClO3	1000415-09-1	43





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111723\
Data File : BP018251.D
Acq On : 18 Nov 2023 09:17
Operator : MA/JU
Sample : 05370-03
Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCDP9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
Quant Title : SVOA CALIBRATION

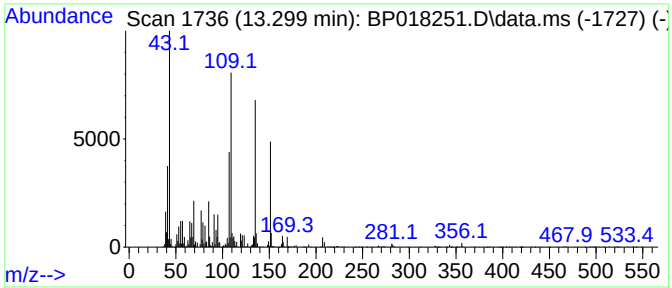
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

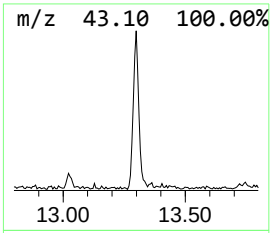
****

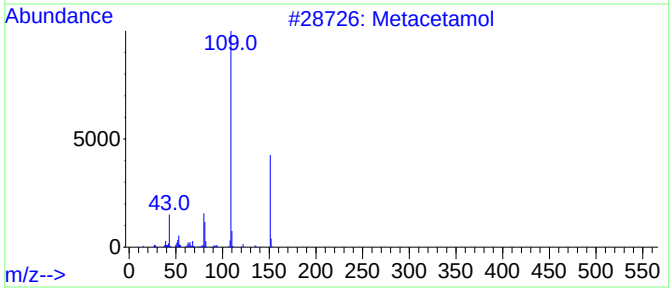
Peak Number 3 Metacetamol Concentration Rank 11

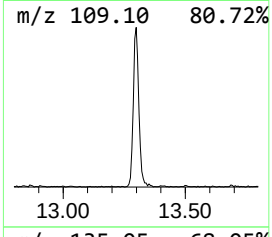
R.T.	EstConc	Area	Relative to ISTD	R.T.
13.299	2.05 ng/ul	82705	Acenaphthene-d10	14.457

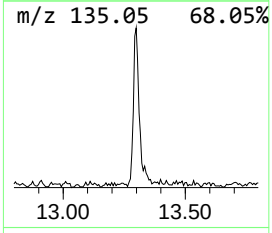
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Metacetamol	151	C8H9NO2	000621-42-1	50
2	Pyridine, 3-butyl-, 1-oxide	151	C9H13NO	031396-33-5	46
3	m-Methoxybenzamide	151	C8H9NO2	005813-86-5	45
4	Metacetamol	151	C8H9NO2	000621-42-1	43
5	Metacetamol	151	C8H9NO2	000621-42-1	43

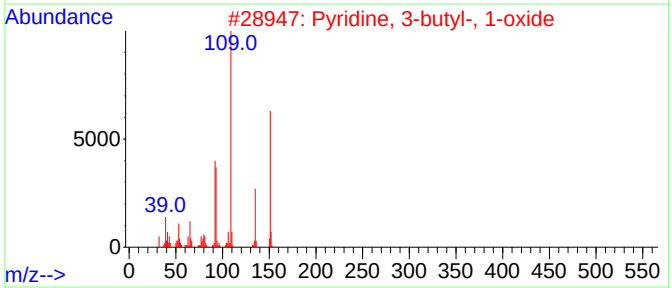


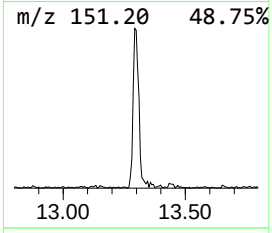


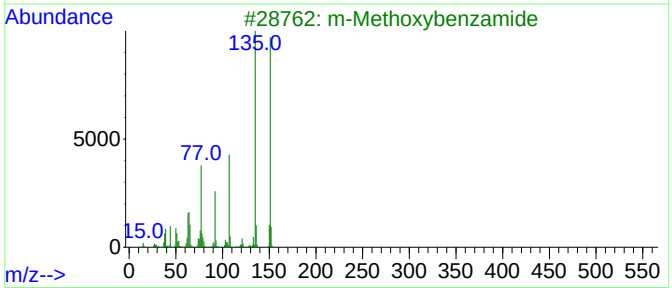


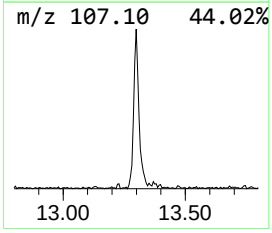












Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111723\
Data File : BP018251.D
Acq On : 18 Nov 2023 09:17
Operator : MA/JU
Sample : 05370-03
Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCDP9

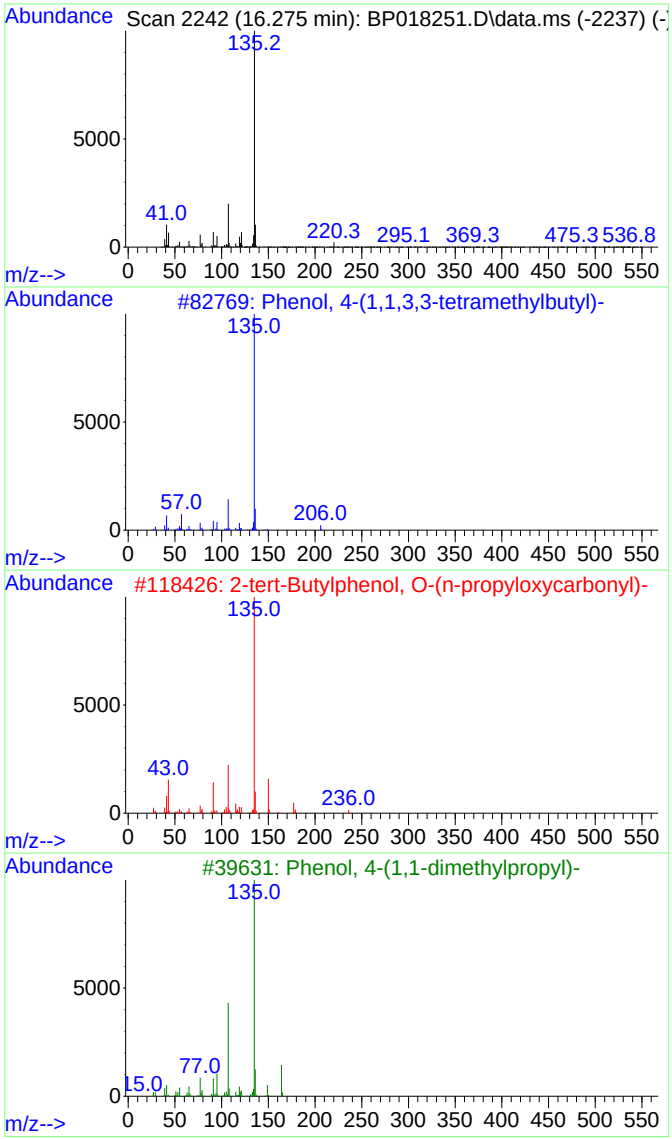
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
Quant Title : SVOA CALIBRATION

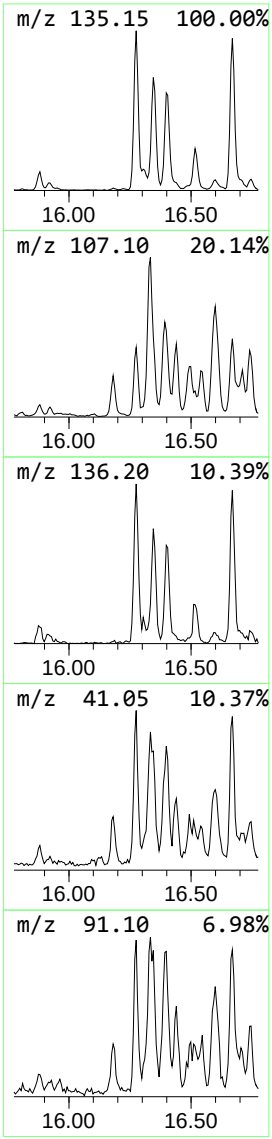
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Phenol, 4-(1,1,3,3-tetramet... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.275	4.05 ng/ul	212727	Phenanthrene-d10	17.245

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 4-(1,1,3,3-tetramethylbu...	206	C14H22O	000140-66-9	86
2	2-tert-Butylphenol, O-(n-propylo...	236	C14H20O3	1000487-37-9	78
3	Phenol, 4-(1,1-dimethylpropyl)-	164	C11H16O	000080-46-6	78
4	Phenol, 2-(1,1,3,3-tetramethylbu...	206	C14H22O	003884-95-5	78
5	Phenol, m-tert-butyl-	150	C10H14O	000585-34-2	78





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Data File : BP018251.D
Acq On : 18 Nov 2023 09:17
Operator : MA/JU
Sample : 05370-03
Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCDP9

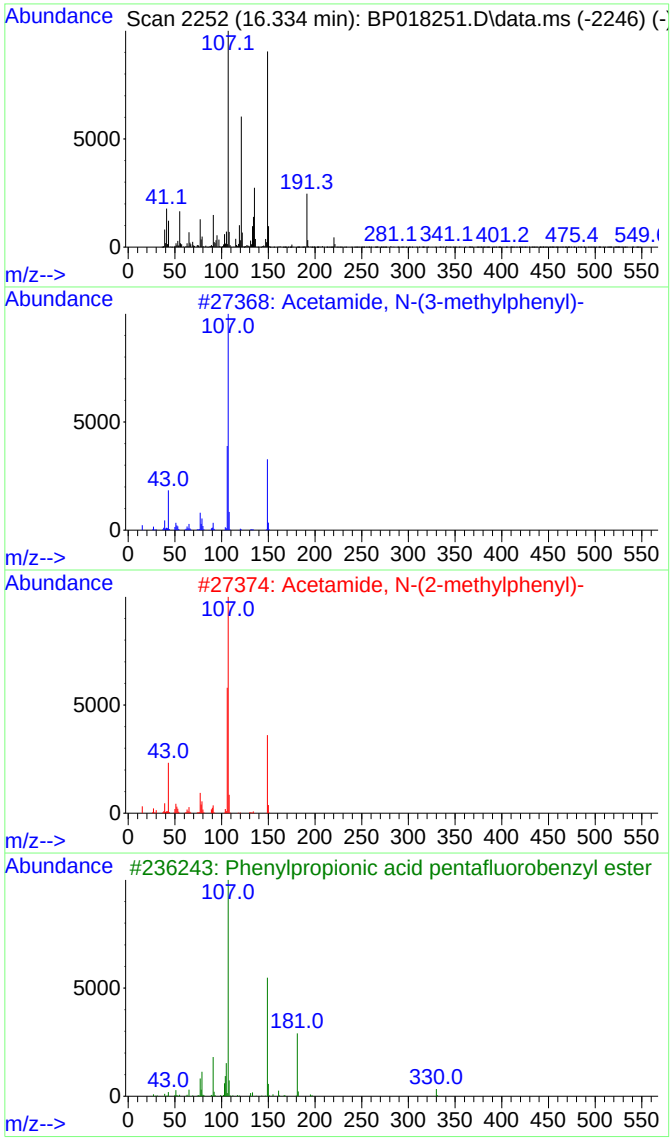
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
Quant Title : SVOA CALIBRATION

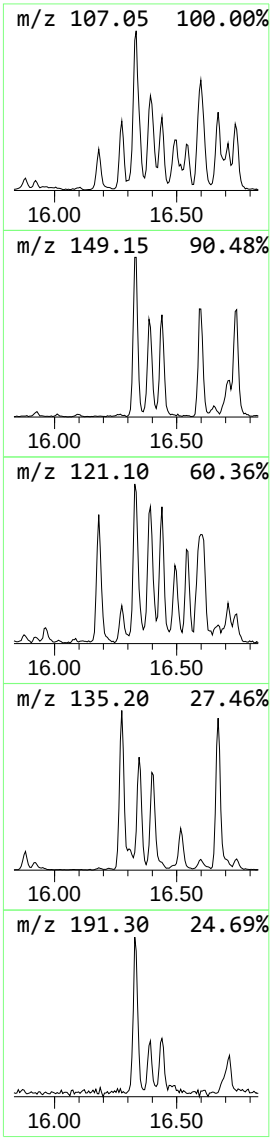
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Acetamide, N-(3-methylphenyl)- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.334	7.50 ng/ul	393942	Phenanthrene-d10	17.245

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Acetamide, N-(3-methylphenyl)-	149	C9H11NO	000537-92-8	52
2	Acetamide, N-(2-methylphenyl)-	149	C9H11NO	000120-66-1	50
3	Phenylpropionic acid pentafluoro...	330	C16H11F5O2	062434-95-1	50
4	N-(Chroman-5-yl)acetamide	191	C11H13NO2	1000306-69-8	45
5	Acetamide, N-(4-methylphenyl)-	149	C9H11NO	000103-89-9	43





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Data File : BP018251.D
Acq On : 18 Nov 2023 09:17
Operator : MA/JU
Sample : 05370-03
Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCDP9

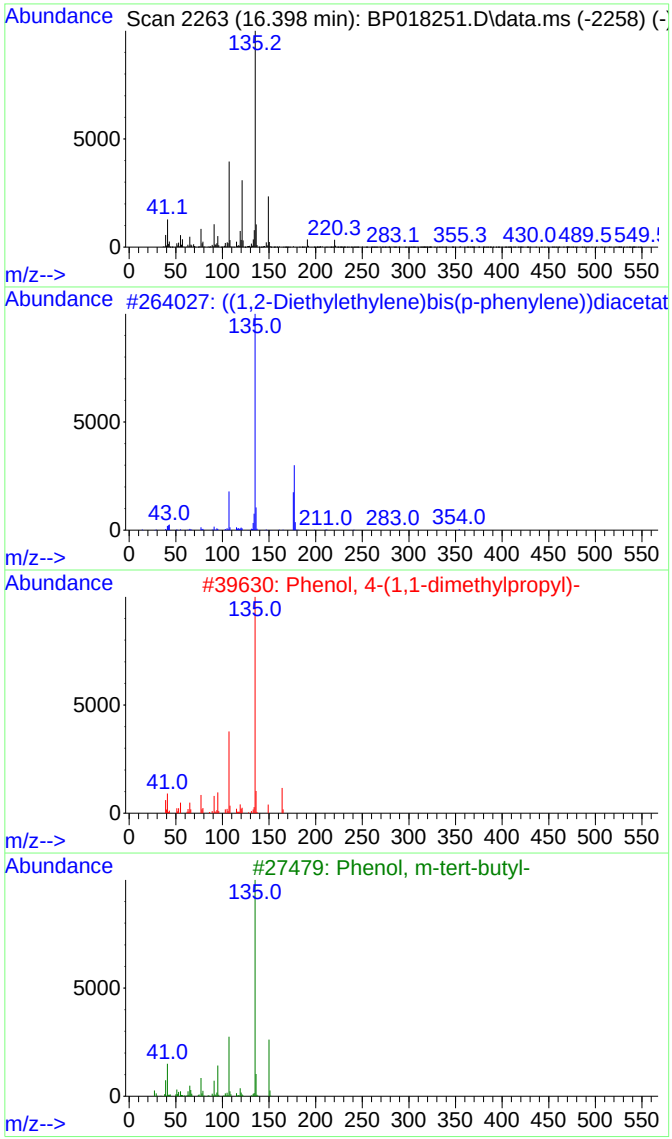
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
Quant Title : SVOA CALIBRATION

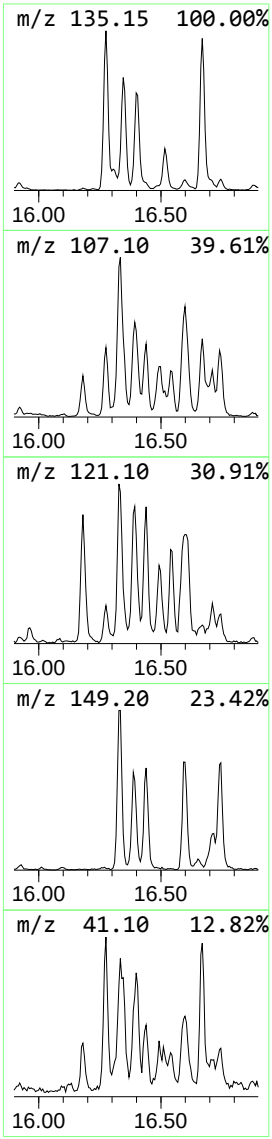
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 ((1,2-Diethylethylene)bis(p... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.398	5.29 ng/ul	277851	Phenanthrene-d10	17.245

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	((1,2-Diethylethylene)bis(p-phen...	354	C22H26O4	004547-76-6	74
2	Phenol, 4-(1,1-dimethylpropyl)-	164	C11H16O	000080-46-6	64
3	Phenol, m-tert-butyl-	150	C10H14O	000585-34-2	64
4	((1,2-Diethylethylene)bis(p-phen...	354	C22H26O4	004547-76-6	64
5	4-(7-Methyloctyl)phenol	220	C15H24O	024518-48-7	60





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111723\
Data File : BP018251.D
Acq On : 18 Nov 2023 09:17
Operator : MA/JU
Sample : 05370-03
Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCDP9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
Quant Title : SVOA CALIBRATION

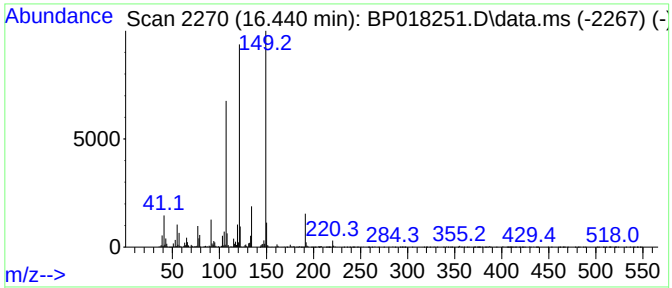
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

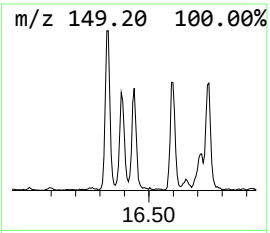
****

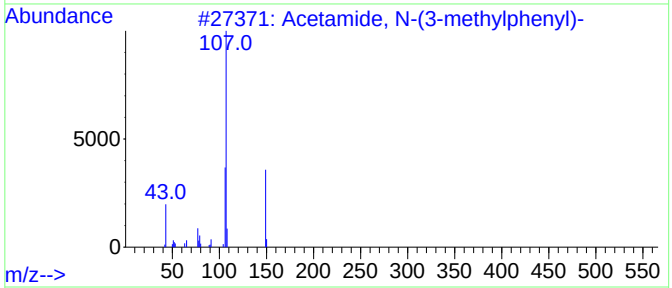
Peak Number 7 unknown-01 Concentration Rank 8

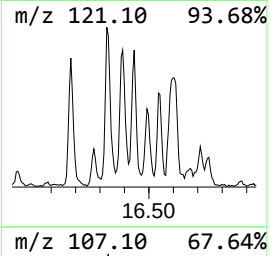
R.T.	EstConc	Area	Relative to ISTD	R.T.
16.440	2.64 ng/ul	138490	Phenanthrene-d10	17.245

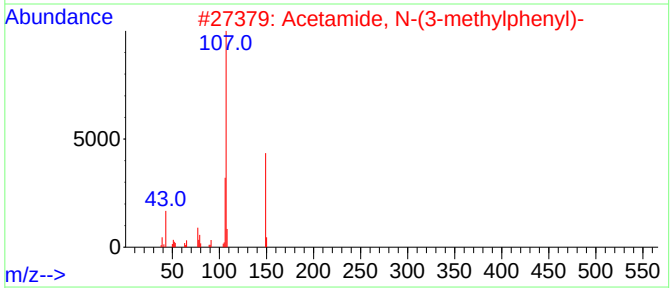
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Acetamide, N-(3-methylphenyl)-	149	C9H11NO	000537-92-8	43
2	Acetamide, N-(3-methylphenyl)-	149	C9H11NO	000537-92-8	43
3	9,10-Dibromo-(+)-camphor	308	C10H14Br2O	085706-53-2	43
4	2-Methyl-4-hydroxybenzoxazole	149	C8H7NO2	051110-60-2	38
5	3-Mercaptoindole, S-acetyl-	191	C10H9NOS	1081542-19-9	38

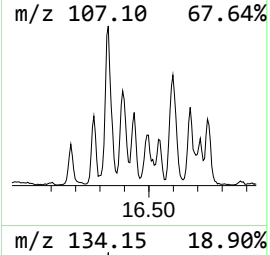


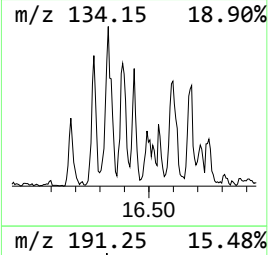


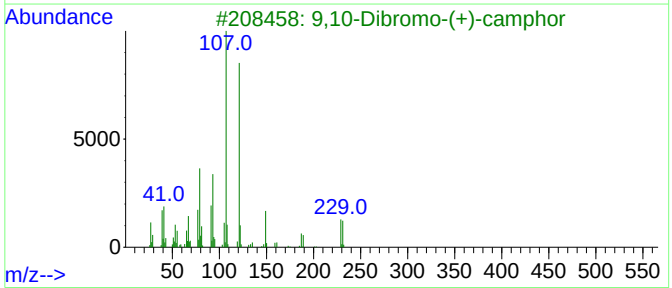


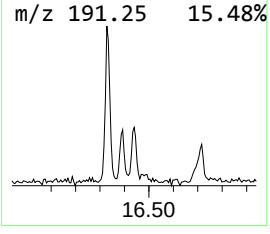












Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111723\
Data File : BP018251.D
Acq On : 18 Nov 2023 09:17
Operator : MA/JU
Sample : 05370-03
Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCDP9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
Quant Title : SVOA CALIBRATION

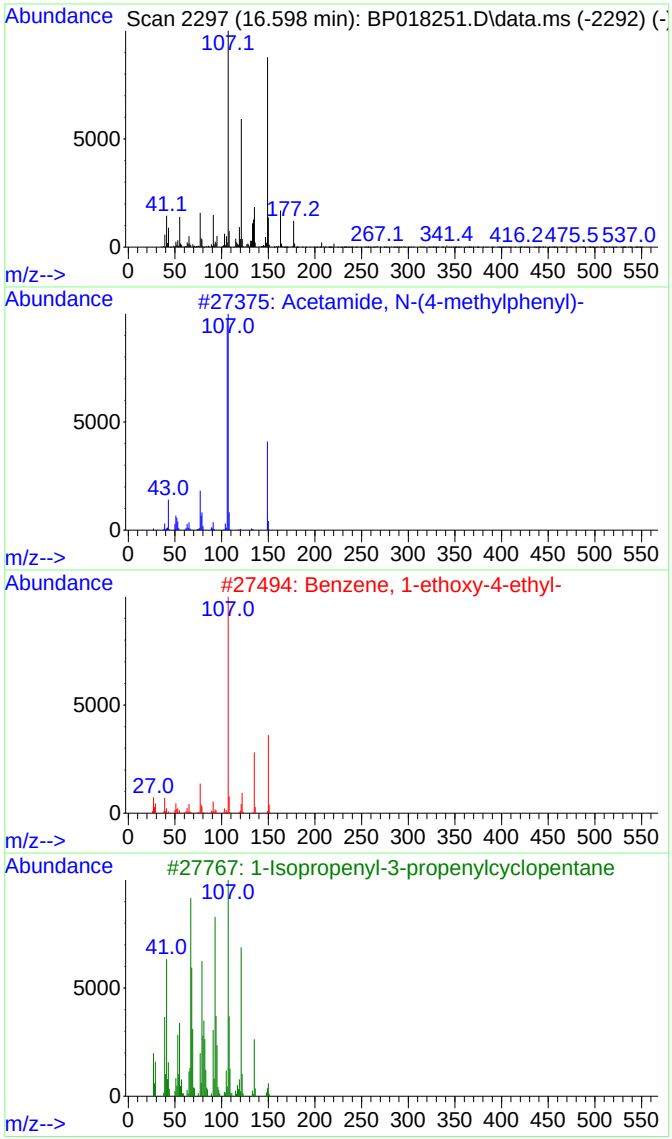
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

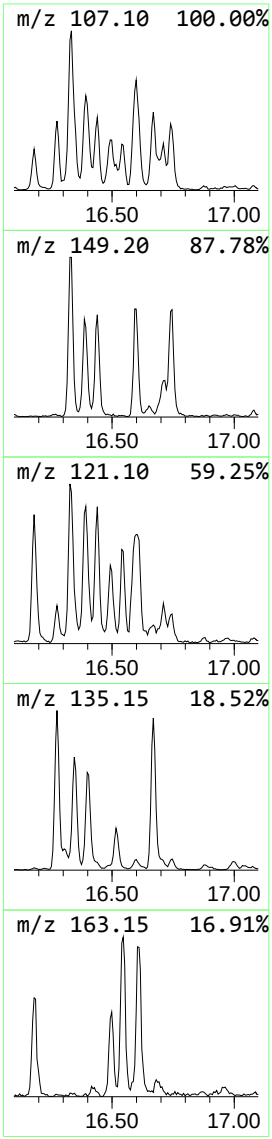
****

Peak Number 8 unknown-02 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.598	4.54 ng/ul	238654	Phenanthrene-d10	17.245

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Acetamide, N-(4-methylphenyl)-	149	C9H11NO	000103-89-9	43
2	Benzene, 1-ethoxy-4-ethyl-	150	C10H14O	001585-06-4	30
3	1-Isopropenyl-3-propenylcyclopen...	150	C11H18	1000192-81-2	27
4	ethanone, 1-(4-methoxy-2-methylp...	240	C16H16O2	1000401-19-7	27
5	Ethanol, 2-[(4-aminophenyl)ethyl...	180	C10H16N2O	000092-65-9	27





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111723\
Data File : BP018251.D
Acq On : 18 Nov 2023 09:17
Operator : MA/JU
Sample : 05370-03
Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCDP9

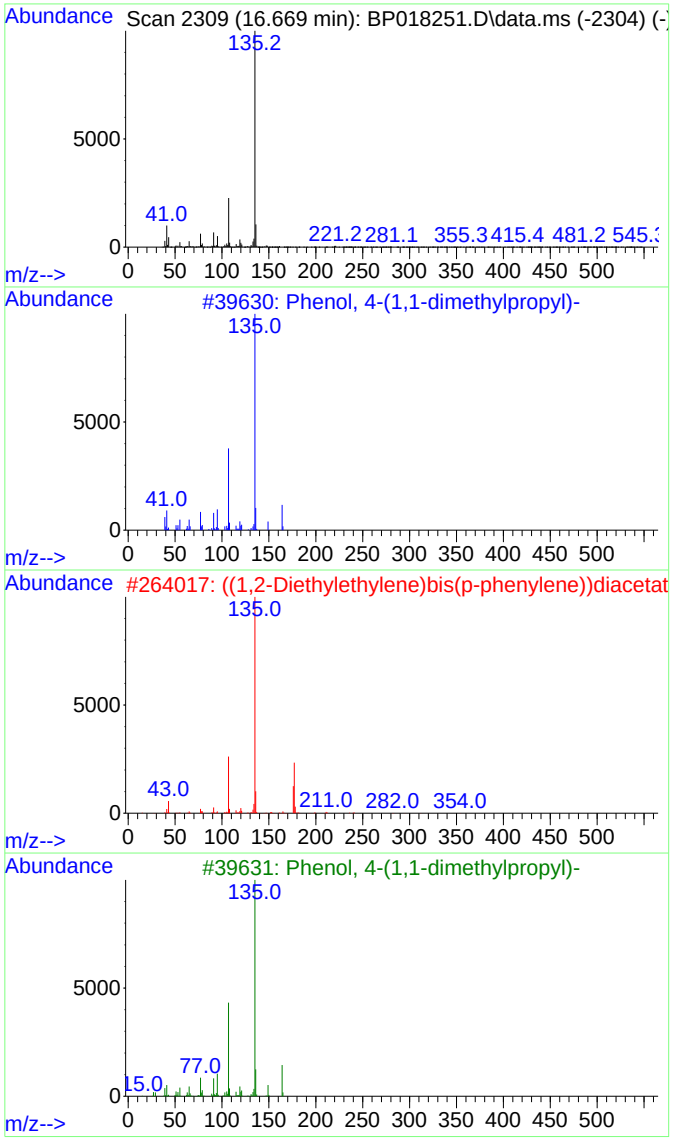
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
Quant Title : SVOA CALIBRATION

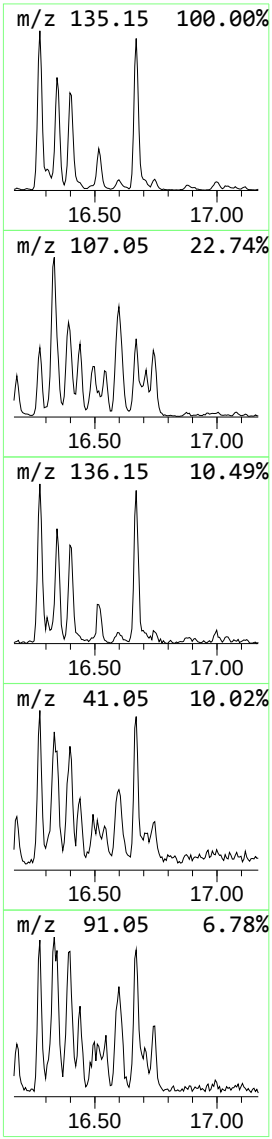
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 Phenol, 4-(1,1-dimethylprop... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.669	4.24 ng/ul	222651	Phenanthrene-d10	17.245

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 4-(1,1-dimethylpropyl)-	164	C11H16O	000080-46-6	83
2	((1,2-Diethylethylene)bis(p-phen...	354	C22H26O4	004547-76-6	78
3	Phenol, 4-(1,1-dimethylpropyl)-	164	C11H16O	000080-46-6	78
4	Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	78
5	4-tert-Butylphenol, O-isopropylo...	236	C14H20O3	1201410-82-3	78





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111723\
Data File : BP018251.D
Acq On : 18 Nov 2023 09:17
Operator : MA/JU
Sample : 05370-03
Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCDP9

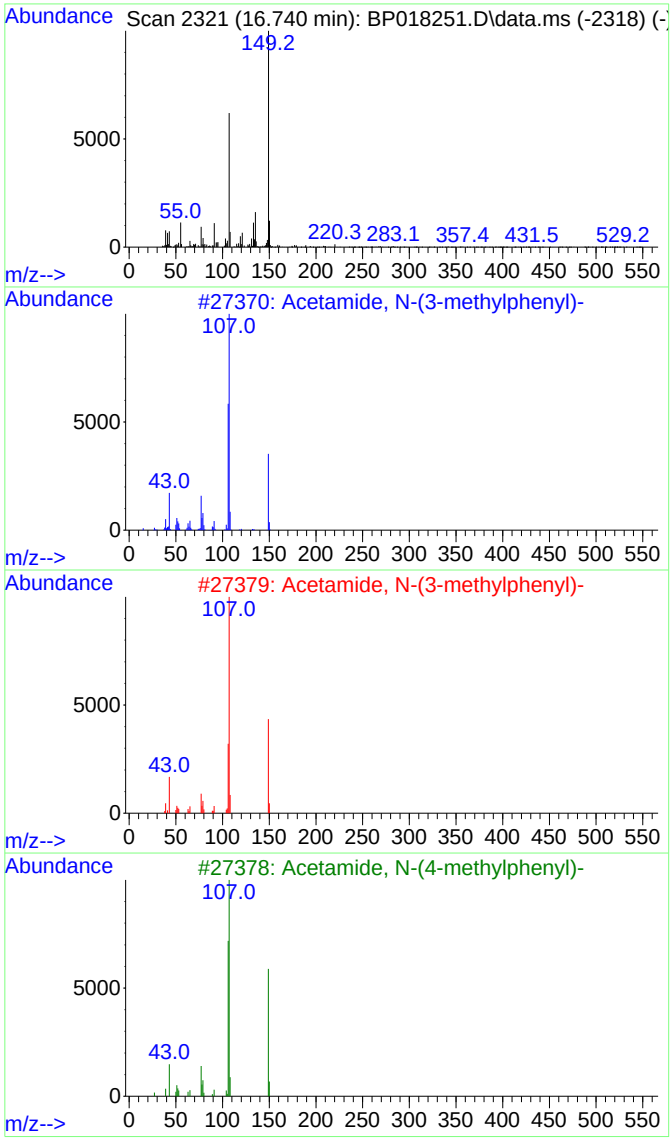
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
Quant Title : SVOA CALIBRATION

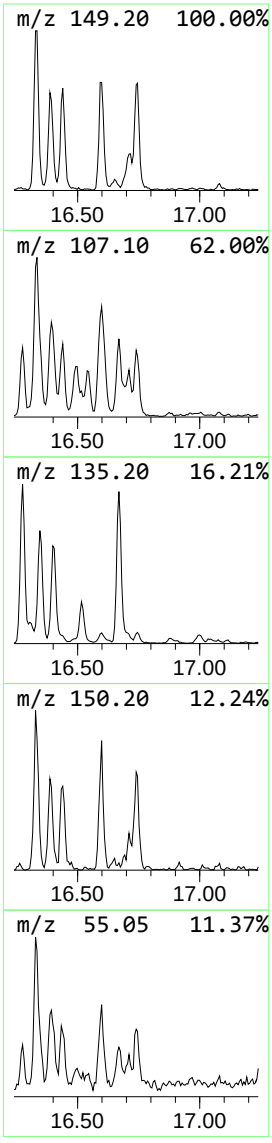
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 Acetamide, N-(4-methylphenyl)- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.740	2.16 ng/ul	113334	Phenanthrene-d10	17.245

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Acetamide, N-(3-methylphenyl)-	149	C9H11NO	000537-92-8	64
2	Acetamide, N-(3-methylphenyl)-	149	C9H11NO	000537-92-8	59
3	Acetamide, N-(4-methylphenyl)-	149	C9H11NO	000103-89-9	59
4	Acetamide, N-(2-methylphenyl)-	149	C9H11NO	000120-66-1	59
5	Ethanone, 2-(1-methylethoxy)-1,2...	254	C17H18O2	006652-28-4	59





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111723\
Data File : BP018251.D
Acq On : 18 Nov 2023 09:17
Operator : MA/JU
Sample : 05370-03
Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCDP9

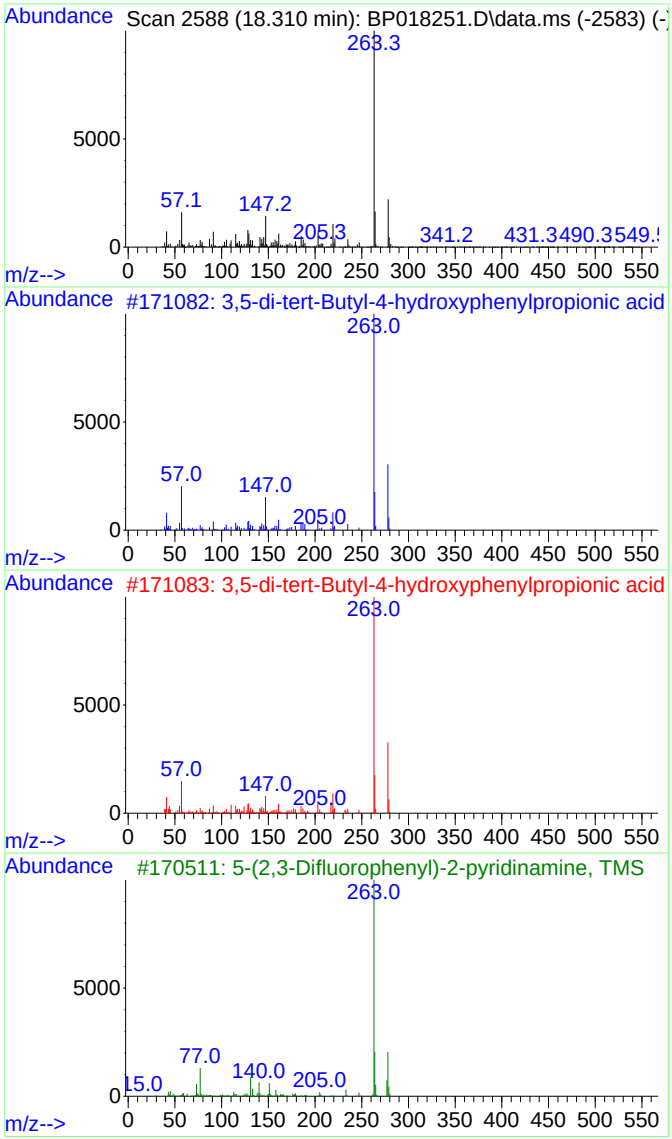
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
Quant Title : SVOA CALIBRATION

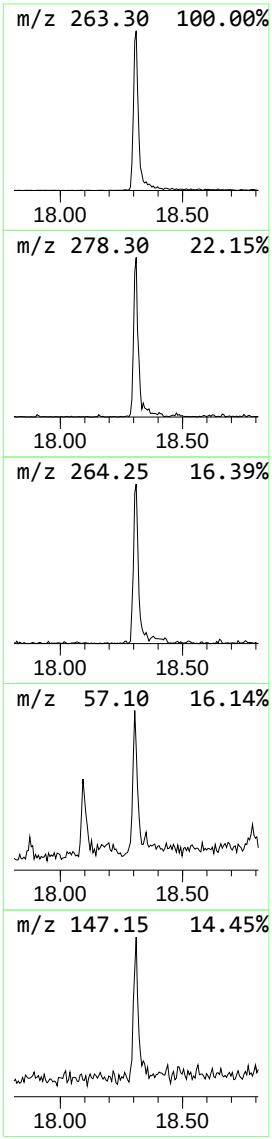
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 11 3,5-di-tert-Butyl-4-hydroxy... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.310	2.68 ng/ul	140645	Phenanthrene-d10	17.245

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3,5-di-tert-Butyl-4-hydroxypheny...	278	C17H26O3	020170-32-5	96
2	3,5-di-tert-Butyl-4-hydroxypheny...	278	C17H26O3	020170-32-5	94
3	5-(2,3-Difluorophenyl)-2-pyridin...	278	C14H16F2N2Si	1000504-49-4	74
4	6,7-Dimethoxy-4(1H)-quinazolinon...	278	C13H18N2O3Si	1000479-65-5	72
5	2'-Hydroxy-4'-methoxyacetophenon...	278	C11H9ClF2O4	1000447-49-1	64





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111723\
Data File : BP018251.D
Acq On : 18 Nov 2023 09:17
Operator : MA/JU
Sample : 05370-03
Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCDP9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

					--Internal Standard---			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
Cyclopentasilox...	9.152	2.2	ng/ul	51804	1	7.781	480131	20.0
Hydrazine, 1-et...	11.405	2.8	ng/ul	85509	2	10.593	617177	20.0
Metacetamol	13.299	2.0	ng/ul	82705	3	14.457	805598	20.0
Phenol, 4-(1,1,...	16.275	4.0	ng/ul	212727	4	17.245	1050640	20.0
Acetamide, N-(3...	16.334	7.5	ng/ul	393942	4	17.245	1050640	20.0
((1,2-Diethylet...	16.398	5.3	ng/ul	277851	4	17.245	1050640	20.0
unknown-01	16.440	2.6	ng/ul	138490	4	17.245	1050640	20.0
unknown-02	16.598	4.5	ng/ul	238654	4	17.245	1050640	20.0
Phenol, 4-(1,1-...	16.669	4.2	ng/ul	222651	4	17.245	1050640	20.0
Acetamide, N-(4...	16.740	2.2	ng/ul	113334	4	17.245	1050640	20.0
3,5-di-tert-But...	18.310	2.7	ng/ul	140645	4	17.245	1050640	20.0