LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071922\
Data File : BP011035.D
Acq On : 19 Jul 2022 16:41
Operator : CG/JU
Sample : N3753-08
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GBHP9

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-BP071422.M
Title : SVOA CALIBRATION

Signal : TIC: BP011035.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.181	12	16	26	rVB	97819	143877	0.82%	0.119%
2	3.599	84	87	104	rVB	324014	527867	2.99%	0.438%
3	3.805	120	122	128	rVB	41163	59358	0.34%	0.049%
4	3.899	135	138	144	rVB	42468	55768	0.32%	0.046%
5	4.264	197	200	205	rVB3	39905	52501	0.30%	0.044%
6	6.858	630	641	642	rBV	305475	433336	2.46%	0.359%
7	6.881	642	645	653	rVB	442062	673848	3.82%	0.559%
8	7.022	662	669	680	rBV	1257625	1946356	11.03%	1.614%
9	7.211	693	701	712	rBV	1176066	1785570	10.12%	1.481%
10	7.675	771	780	787	rBV	1313268	2011650	11.40%	1.669%
11	7.752	787	793	800	rVV	123304	198270	1.12%	0.164%
12	7.928	816	823	831	rBV	155527	279617	1.58%	0.232%
13	8.393	893	902	913	rBV	850703	1336387	7.57%	1.109%
14	8.840	970	978	987	rBV	1352082	2179510	12.35%	1.808%
15	8.987	994	1003	1014	rBV3	273271	620824	3.52%	0.515%
16	9.257	1043	1049	1055	rVB	69726	109287	0.62%	0.091%
17	9.557	1093	1100	1108	rBV	961136	1537557	8.71%	1.275%
18	9.716	1121	1127	1136	rVB2	27511	56088	0.32%	0.047%
19	9.804	1138	1142	1147	rBV3	30204	49867	0.28%	0.041%
20	10.093	1184	1191	1202	rBV	1489946	2438252	13.82%	2.022%
21	10.469	1248	1255	1267	rVB	1700206	2812716	15.94%	2.333%
22	10.616	1269	1280	1290	rBV	976587	1762478	9.99%	1.462%
23	10.687	1290	1292	1302	rVB6	31894	68698	0.39%	0.057%
24	10.840	1311	1318	1345	rBV	2602197	4724192	26.77%	3.919%
25	11.022	1345	1349	1355	rVV	30223	56416	0.32%	0.047%
26	11.087	1355	1360	1363	rVV	33057	55718	0.32%	0.046%
27	11.122	1364	1366	1372	rVB2	27330	41460	0.23%	0.034%
28	11.228	1377	1384	1390	rBV	162080	290190	1.64%	0.241%
29	11.287	1390	1394	1403	rVB2	87642	147767	0.84%	0.123%
30	11.446	1414	1421	1430	rBV3	44015	99244	0.56%	0.082%
31	11.757	1468	1474	1482	rVB3	41771	70430	0.40%	0.058%
32	12.069	1520	1527	1532	rVB3	40162	74499	0.42%	0.062%
33	12.622	1616	1621	1627	rBV5	27813	42348	0.24%	0.035%
34	12.946	1670	1676	1682	rVB2	31917	49883	0.28%	0.041%
35	13.181	1709	1716	1722	rBV	328735	470611	2.67%	0.390%
36	13.245	1722	1727	1734	rVV2	117185	216699	1.23%	0.180%

LSC Area Percent Report

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Acq On : 19 Jul 2022 16:41
Operator : CG/JU
Sample : N3753-08
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GBHP9

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-BP071422.M
Title : SVOA CALIBRATION

37	13.310	1734	1738	1744	rVB3	44872	68560	0.39%	0.057%
38	13.757	1807	1814	1820	rBV	2950461	4035078	22.86%	3.347%
39	13.804	1820	1822	1827	rVV	36711	49372	0.28%	0.041%
40	13.875	1827	1834	1844	rVB7	47429	108913	0.62%	0.090%
41	14.028	1850	1860	1871	rBV	3291333	4954465	28.07%	4.110%
42	14.210	1885	1891	1901	rBV2	103311	326150	1.85%	0.271%
43	14.357	1906	1916	1925	rVV2	11112126	17648385	100.00%	14.639%
44	14.504	1936	1941	1946	rBV	513051	664987	3.77%	0.552%
45	14.557	1946	1950	1960	rVB	155853	267773	1.52%	0.222%
46	14.775	1983	1987	1991	rVV2	113790	150041	0.85%	0.124%
47	14.816	1991	1994	2000	rVV3	38255	55969	0.32%	0.046%
48	14.945	2012	2016	2019	rBV3	41856	57433	0.33%	0.048%
49	14.998	2019	2025	2031	rVB3	38398	74360	0.42%	0.062%
50	15.098	2038	2042	2046	rVB	40957	56579	0.32%	0.047%
51	15.157	2048	2052	2054	rBV	51068	70115	0.40%	0.058%
52	15.181	2054	2056	2060	rVB	51401	55750	0.32%	0.046%
53	15.234	2060	2065	2069	rBV	97901	126570	0.72%	0.105%
54	15.339	2076	2083	2097	rVB	4265354	5920943	33.55%	4.911%
55	15.463	2097	2104	2120	rBV	1155310	1746740	9.90%	1.449%
56	15.581	2120	2124	2130	rVB	55425	81743	0.46%	0.068%
57	15.716	2143	2147	2151	rVB2	31335	41426	0.23%	0.034%
58	15.910	2175	2180	2185	rBV3	99800	183107	1.04%	0.152%
59	15.951	2185	2187	2190	rVV4	45011	56955	0.32%	0.047%
60	16.004	2190	2196	2206	rVB	3667127	4855465	27.51%	4.028%
61	16.163	2220	2223	2233	rVB	33034	62919	0.36%	0.052%
62	16.251	2234	2238	2245	rVB6	26619	41062	0.23%	0.034%
63	16.375	2254	2259	2267	rBV3	24426	50261	0.28%	0.042%
64	16.516	2280	2283	2289	rVB4	41996	56118	0.32%	0.047%
65	17.039	2368	2372	2376	rVV	72323	93270	0.53%	0.077%
66	17.104	2377	2383	2390	rVV	2708996	3708759	21.01%	3.076%
67	17.204	2394	2400	2413	rVV2	4694532	6673678	37.81%	5.536%
68	17.434	2431	2439	2445	rBV3	115107	234339	1.33%	0.194%
69	17.516	2445	2453	2459	rBV4	96744	166151	0.94%	0.138%
70	17.616	2467	2470	2475	rVB4	37435	50777	0.29%	0.042%
71	17.733	2486	2490	2495	rBV2	87198	116605	0.66%	0.097%
72	17.792	2495	2500	2512	rVB2	377944	576237	3.27%	0.478%
73	17.910	2513	2520	2528	rBV9	31985	88133	0.50%	0.073%
74	18.022	2534	2539	2544	rBV	192855	268356	1.52%	0.223%
75	18.081	2545	2549	2559	rVB2	201838	353124	2.00%	0.293%
76	18.304	2583	2587	2596	rBV4	48724	83187	0.47%	0.069%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071922\
Data File : BP011035.D
Acq On : 19 Jul 2022 16:41
Operator : CG/JU
Sample : N3753-08
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GBHP9

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_P\Methods\SFAM-EPA-BP071422.M
Title : SVOA CALIBRATION

77	18.581	2630	2634	2642	rVB2	78597	102330	0.58%	0.085%
78	18.863	2677	2682	2689	rBV	571532	842472	4.77%	0.699%
79	18.945	2692	2696	2704	rVV	262196	339852	1.93%	0.282%
80	19.039	2708	2712	2715	rBV2	71628	91572	0.52%	0.076%
81	19.075	2715	2718	2721	rVV	259024	274872	1.56%	0.228%
82	19.128	2721	2727	2730	rVV2	92329	195787	1.11%	0.162%
83	19.169	2730	2734	2738	rVV	469744	528343	2.99%	0.438%
84	19.204	2738	2740	2743	rVB2	38551	42322	0.24%	0.035%
85	19.263	2748	2750	2758	rVB4	46240	69635	0.39%	0.058%
86	19.416	2773	2776	2779	rBV	56019	55023	0.31%	0.046%
87	19.463	2779	2784	2791	rVB	6416474	8124728	46.04%	6.739%
88	19.616	2806	2810	2817	rBV	183053	244679	1.39%	0.203%
89	19.986	2870	2873	2876	rBV2	85247	113608	0.64%	0.094%
90	20.263	2916	2920	2925	rBV	507992	555740	3.15%	0.461%
91	20.333	2929	2932	2935	rBV	133971	134707	0.76%	0.112%
92	20.386	2935	2941	2950	rBV	2684615	3428445	19.43%	2.844%
93	20.492	2956	2959	2963	rBV2	224371	270055	1.53%	0.224%
94	20.957	3034	3038	3045	rBV	965277	1314557	7.45%	1.090%
95	21.022	3045	3049	3057	rVB	1156592	1316188	7.46%	1.092%
96	21.227	3079	3084	3098	rBV	3481850	4446269	25.19%	3.688%
97	21.769	3173	3176	3181	rVB2	438096	490129	2.78%	0.407%
98	22.339	3268	3273	3285	rBV	1884294	2903328	16.45%	2.408%
99	23.433	3452	3459	3468	rBV2	4444839	8449847	47.88%	7.009%
100	23.580	3478	3484	3495	rVB2	2181716	4462886	25.29%	3.702%

Sum of corrected areas: 120556368

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071922\
Data File : BP011035.D
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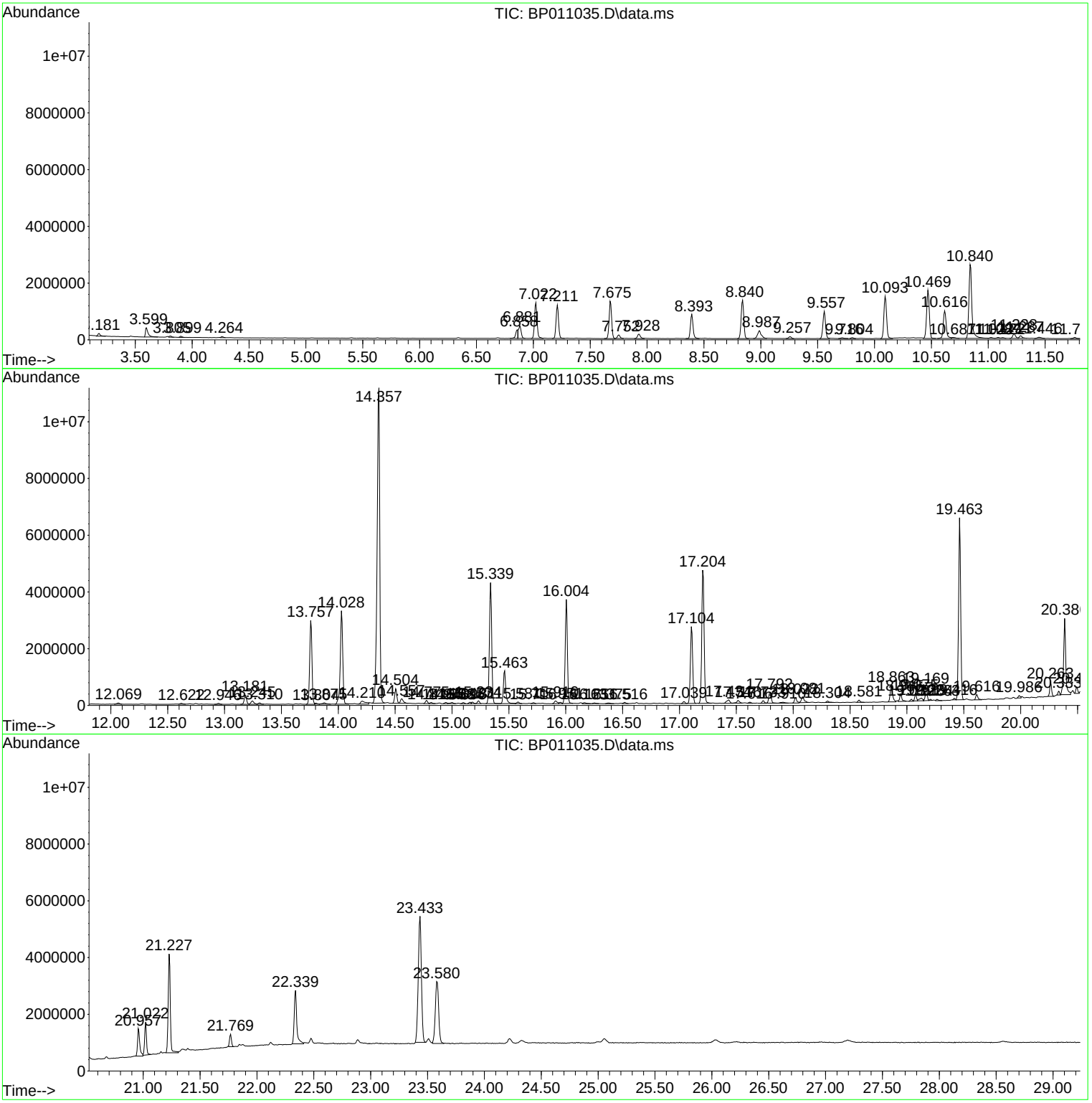
Operator : CG/JU
Sample : N3753-08
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GBHP9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071422.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071922\
Data File : BP011035.D
Acq On : 19 Jul 2022 16:41
Operator : CG/JU
Sample : N3753-08
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GBHP9

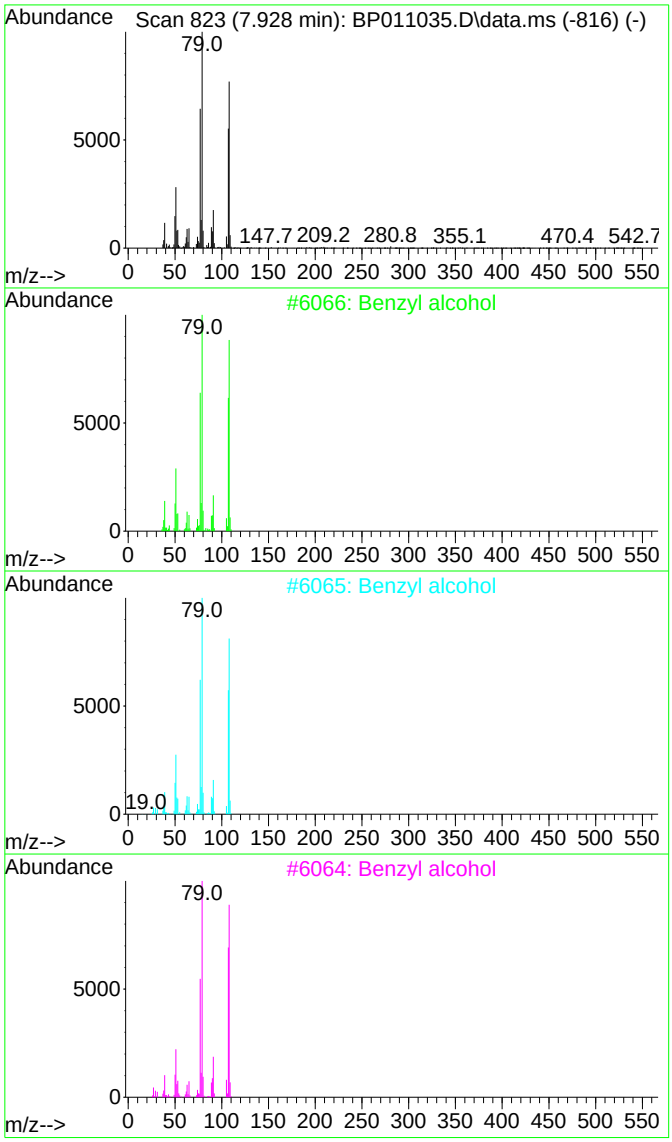
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071422.M
Quant Title : SVOA CALIBRATION

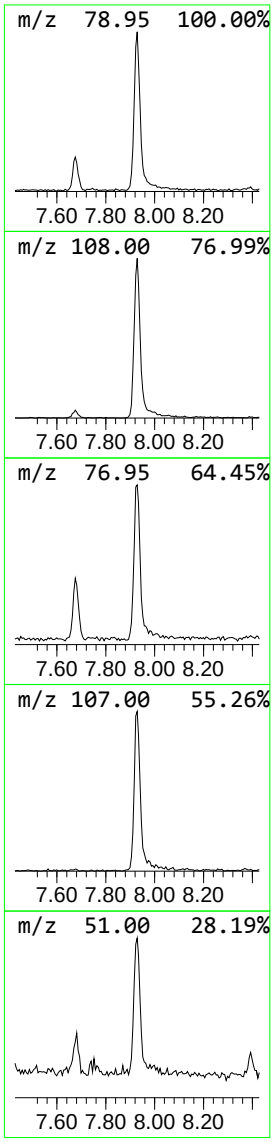
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 Benzyl alcohol Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.928	2.78 ng/ul	279617	1,4-Dichlorobenzene-d4	7.675

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzyl alcohol	108	C7H8O	000100-51-6	98
2	Benzyl alcohol	108	C7H8O	000100-51-6	98
3	Benzyl alcohol	108	C7H8O	000100-51-6	97
4	Benzyl alcohol	108	C7H8O	000100-51-6	95
5	N-Cbz-glycylglycine p-nitropheny...	387	C18H17N3O7	013574-81-7	91





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071922\
Data File : BP011035.D
Acq On : 19 Jul 2022 16:41
Operator : CG/JU
Sample : N3753-08
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GBHP9

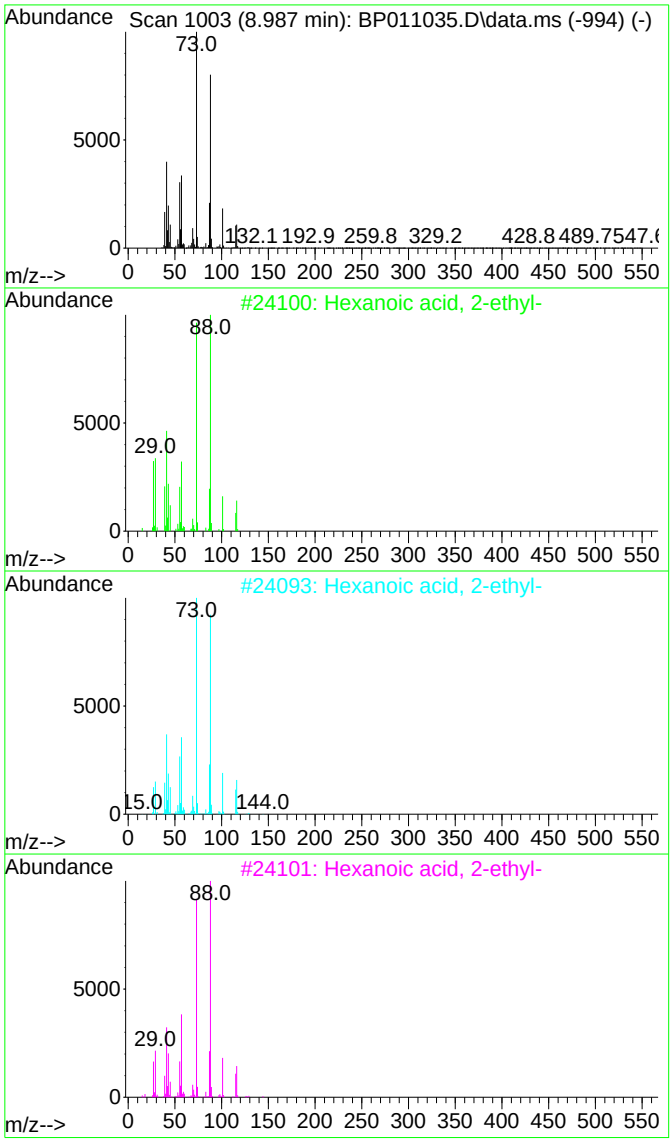
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071422.M
Quant Title : SVOA CALIBRATION

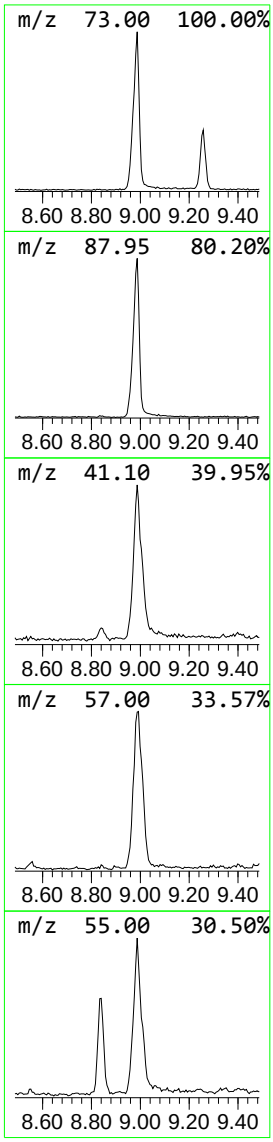
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 Hexanoic acid, 2-ethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.987	6.17 ng/ul	620824	1,4-Dichlorobenzene-d4	7.675

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	90
2	Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	86
3	Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	86
4	Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	78
5	Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	72





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071922\
Data File : BP011035.D
Acq On : 19 Jul 2022 16:41
Operator : CG/JU
Sample : N3753-08
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GBHP9

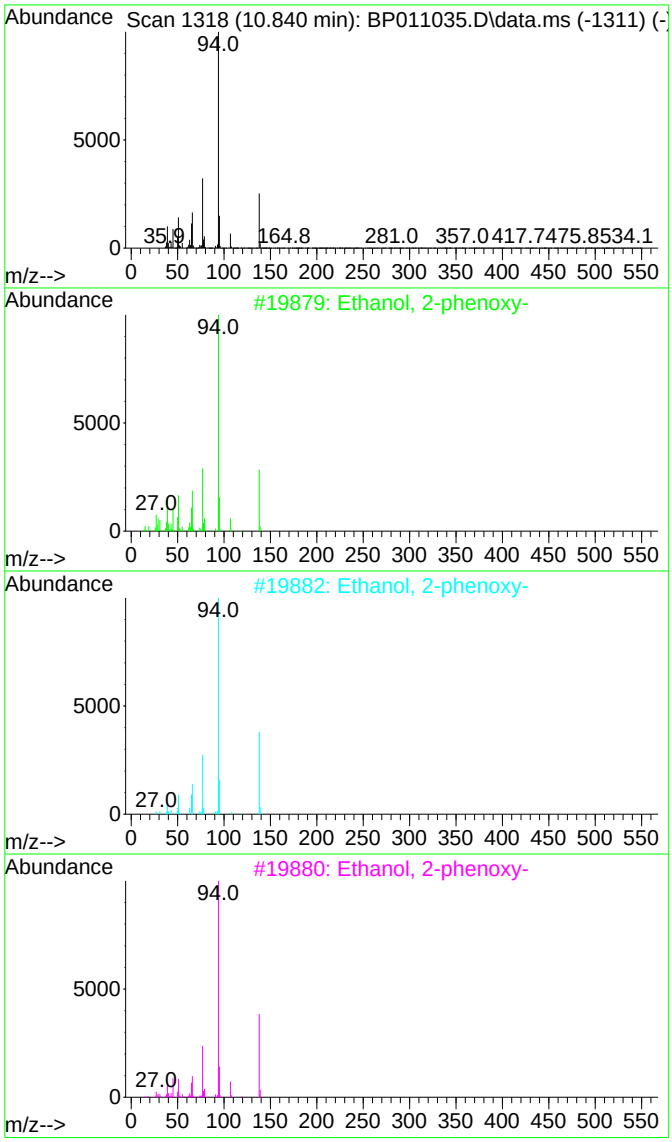
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071422.M
Quant Title : SVOA CALIBRATION

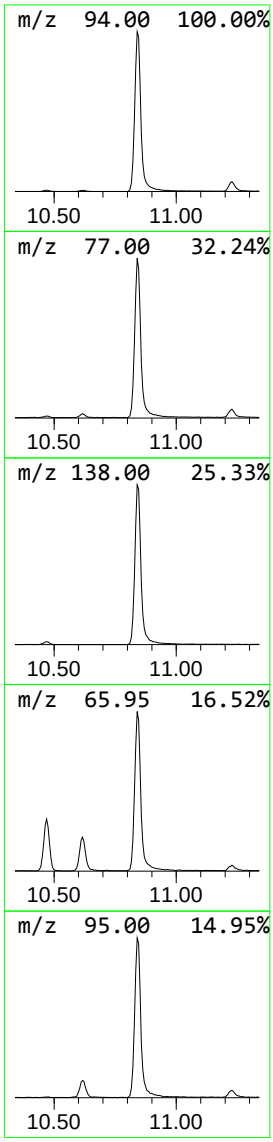
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Ethanol, 2-phenoxy- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.840	33.59 ng/ul	4724190	Naphthalene-d8	10.469

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethanol, 2-phenoxy-	138	C8H10O2	000122-99-6	97
2	Ethanol, 2-phenoxy-	138	C8H10O2	000122-99-6	94
3	Ethanol, 2-phenoxy-	138	C8H10O2	000122-99-6	91
4	Ethanol, 2-phenoxy-	138	C8H10O2	000122-99-6	91
5	Carbonic acid, neopentyl phenyl ...	208	C12H16O3	013183-19-2	64





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071922\
Data File : BP011035.D
Acq On : 19 Jul 2022 16:41
Operator : CG/JU
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Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GBHP9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071422.M
Quant Title : SVOA CALIBRATION

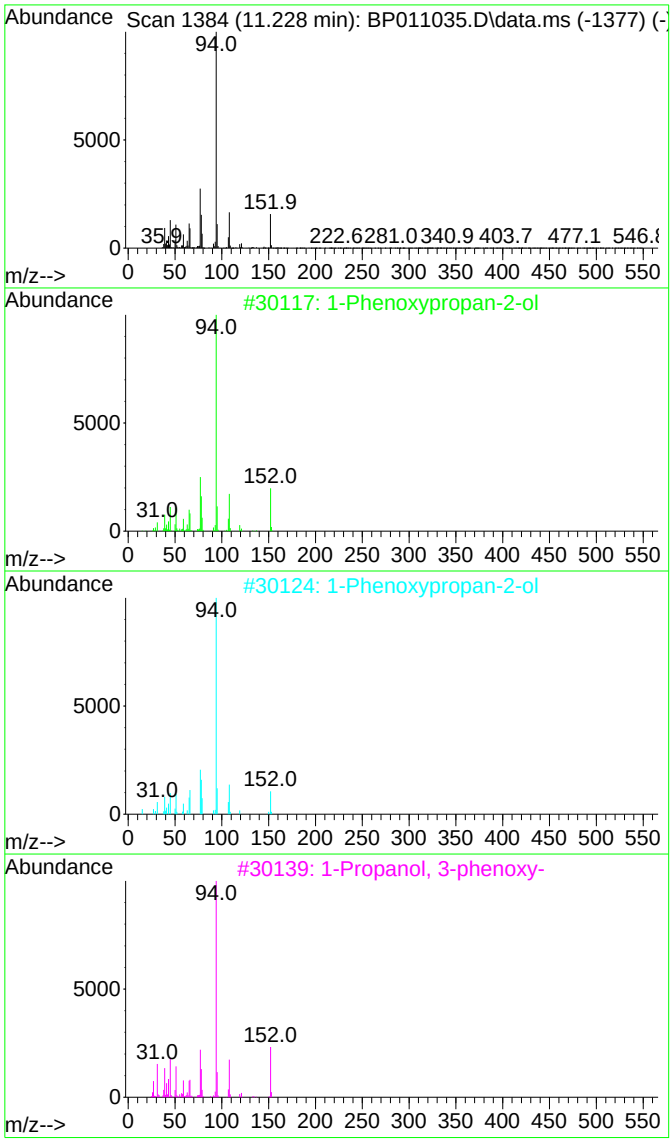
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

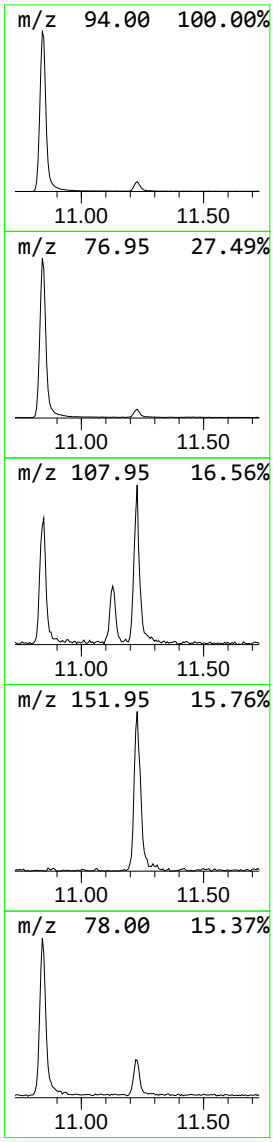
****

Peak Number 4 1-Phenoxypropan-2-ol Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.228	2.06 ng/ul	290190	Naphthalene-d8	10.469

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	96
2	1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	94
3	1-Propanol, 3-phenoxy-	152	C9H12O2	006180-61-6	94
4	1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	91
5	1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	87





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071922\
Data File : BP011035.D
Acq On : 19 Jul 2022 16:41
Operator : CG/JU
Sample : N3753-08
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GBHP9

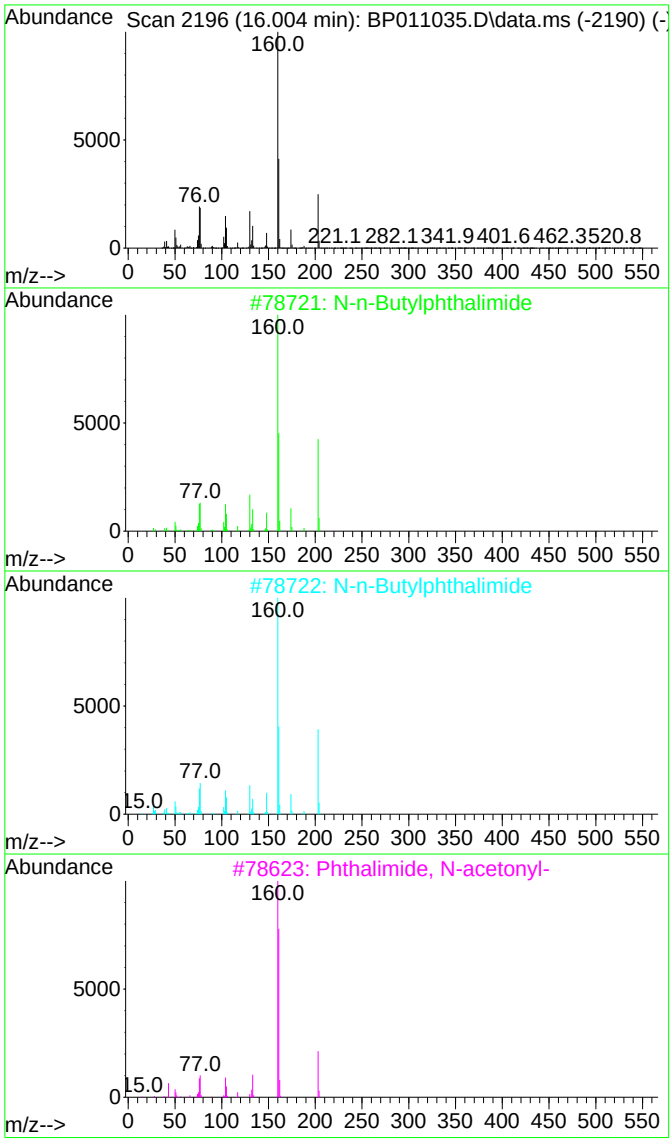
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071422.M
Quant Title : SVOA CALIBRATION

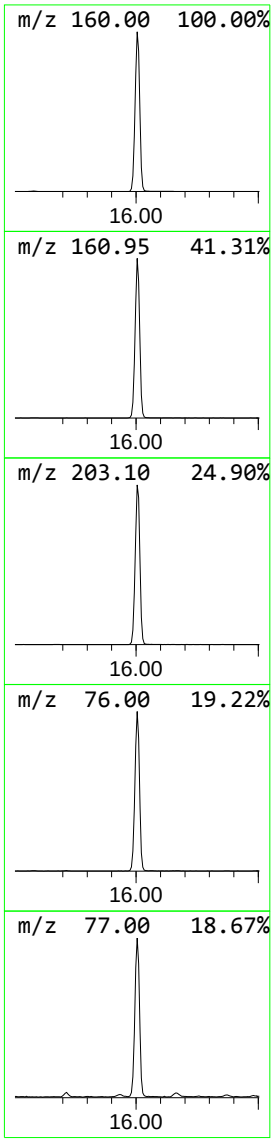
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 N-n-Butylphthalimide Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.004	26.18 ng/ul	4855470	Phenanthrene-d10	17.104

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	N-n-Butylphthalimide	203	C12H13NO2	001515-72-6	98
2	N-n-Butylphthalimide	203	C12H13NO2	001515-72-6	95
3	Phthalimide, N-acetonyl-	203	C11H9NO3	003416-57-7	62
4	(4-Chloro-5-fluoropyridin-2-yl)m...	161	C6H5ClFNO	113209-90-8	59
5	1H-Isoindole-1,3(2H)-dione, 2-(4...	281	C12H12BrNO2	005394-18-3	53





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Data File : BP011035.D
Acq On : 19 Jul 2022 16:41
Operator : CG/JU
Sample : N3753-08
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GBHP9

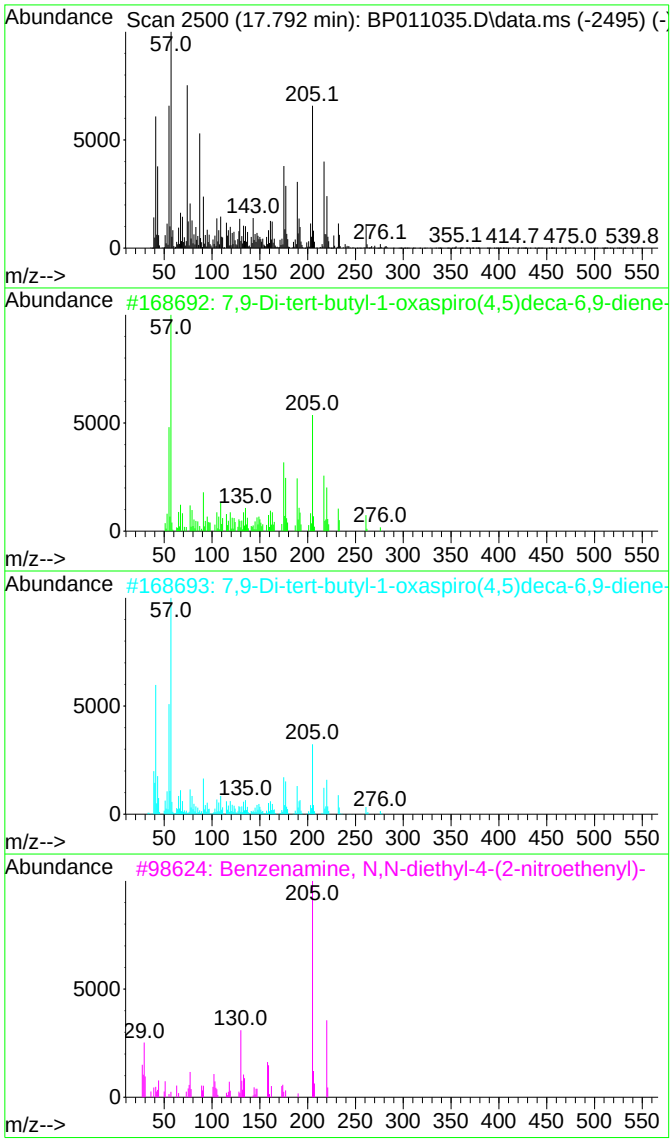
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071422.M
Quant Title : SVOA CALIBRATION

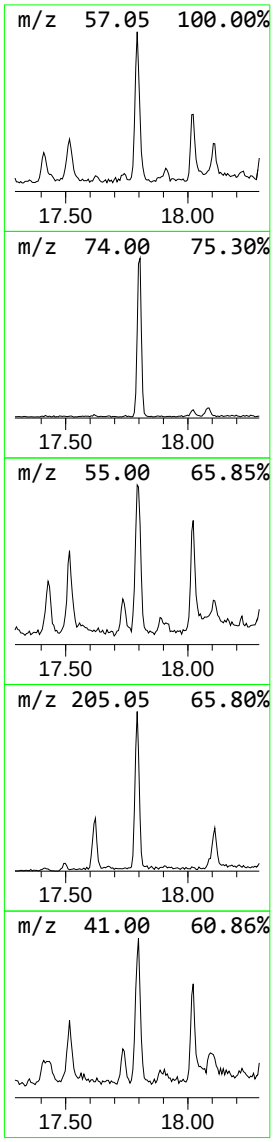
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 7,9-Di-tert-butyl-1-oxaspir... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.792	3.11 ng/ul	576237	Phenanthrene-d10	17.104

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	7,9-Di-tert-butyl-1-oxaspiro(4,5...	276	C17H24O3	082304-66-3	98
2	7,9-Di-tert-butyl-1-oxaspiro(4,5...	276	C17H24O3	082304-66-3	93
3	Benzenamine, N,N-diethyl-4-(2-ni...	220	C12H16N2O2	064462-51-7	44
4	Methyl 8-methyl-decanoate	200	C12H24O2	1010336-49-1	38
5	7,9-Di-tert-butyl-1-oxaspiro(4,5...	276	C17H24O3	082304-66-3	35





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Operator : CG/JU
Sample : N3753-08
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GBHP9

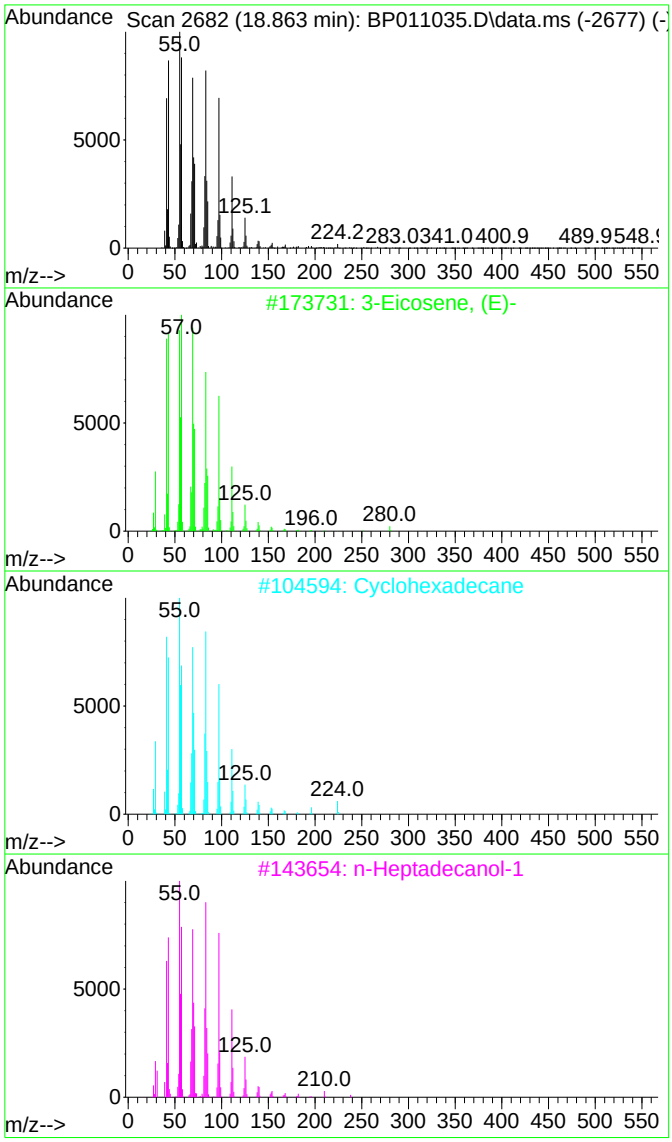
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071422.M
Quant Title : SVOA CALIBRATION

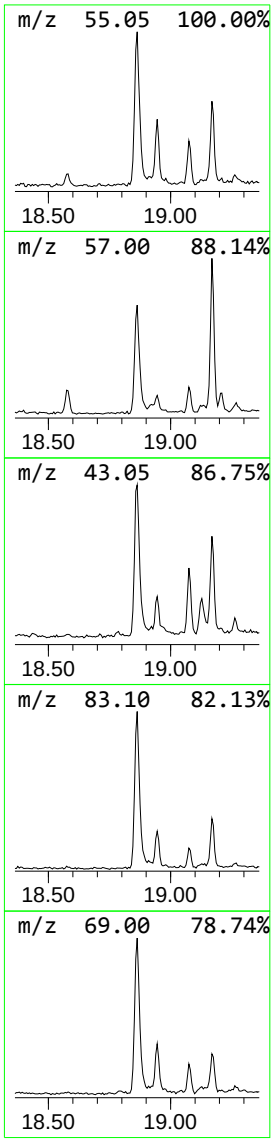
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.863	4.54 ng/ul	842472	Phenanthrene-d10	17.104

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Eicosene, (E)-	280	C20H40	074685-33-9	95
2	Cyclohexadecane	224	C16H32	000295-65-8	94
3	n-Heptadecanol-1	256	C17H36O	001454-85-9	94
4	1-Octadecanol	270	C18H38O	000112-92-5	94
5	5-Eicosene, (E)-	280	C20H40	074685-30-6	93





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Operator : CG/JU
Sample : N3753-08
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GBHP9

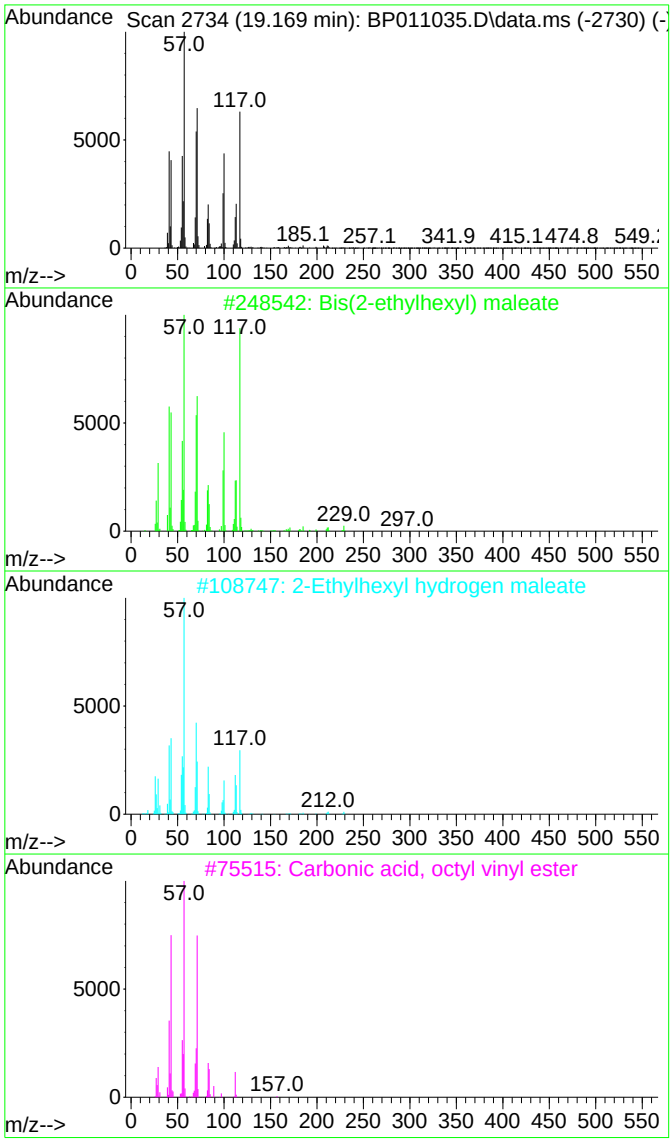
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071422.M
Quant Title : SVOA CALIBRATION

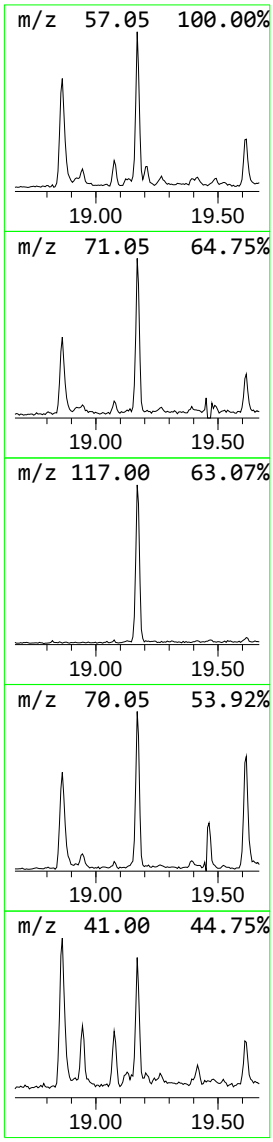
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 Bis(2-ethylhexyl) maleate Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.169	2.38 ng/ul	528343	Chrysene-d12	21.227

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Bis(2-ethylhexyl) maleate	340	C20H36O4	000142-16-5	52
2	2-Ethylhexyl hydrogen maleate	228	C12H20O4	002370-71-0	49
3	Carbonic acid, octyl vinyl ester	200	C11H20O3	1000383-25-5	38
4	1-Decene, 3,4-dimethyl-	168	C12H24	050871-03-9	35
5	Methoxyacetic acid, 2-ethylhexyl...	202	C11H22O3	1000282-41-4	35





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Acq On : 19 Jul 2022 16:41
Operator : CG/JU
Sample : N3753-08
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GBHP9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071422.M
Quant Title : SVOA CALIBRATION

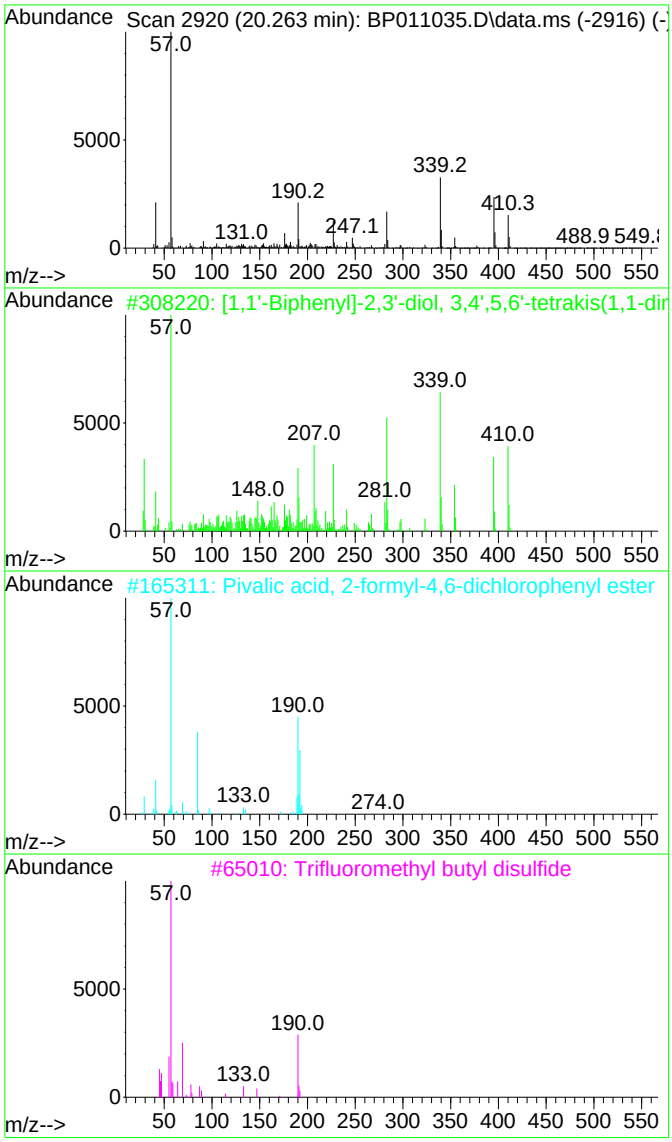
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

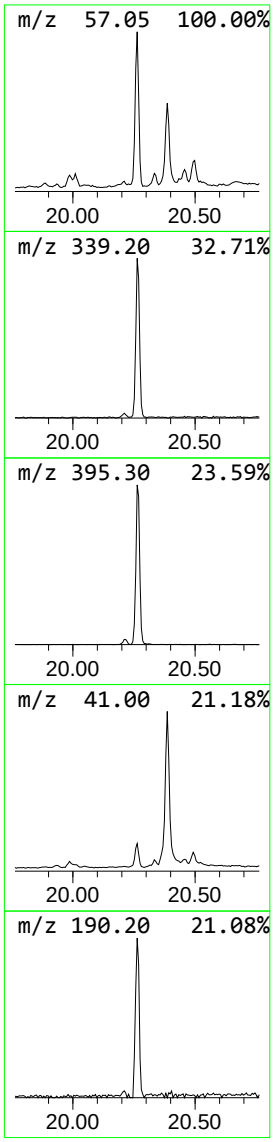
****

Peak Number 9 unknown-01 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.263	2.50 ng/ul	555740	Chrysene-d12	21.227

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	[1,1'-Biphenyl]-2,3'-diol, 3,4',...	410	C28H42O2	1000362-13-2	38
2	Pivalic acid, 2-formyl-4,6-dichl...	274	C12H12Cl2O3	1000330-92-2	9
3	Trifluoromethyl butyl disulfide	190	C5H9F3S2	147618-69-7	9
4	Phosphine sulfide, t-butyl-dichl...	190	C4H9Cl2PS	021187-18-8	9
5	3H-Cycloprop(1,2)androsta-1,4,6-...	395	C24H29NO4	075857-75-9	7





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Operator : CG/JU
Sample : N3753-08
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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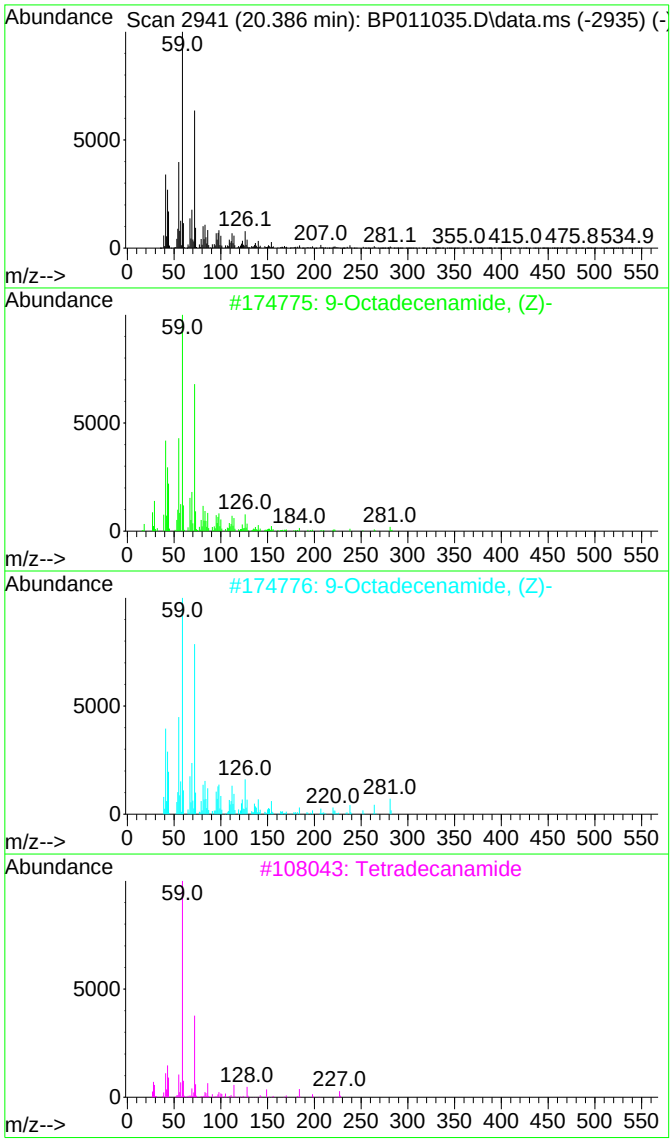
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071422.M
Quant Title : SVOA CALIBRATION

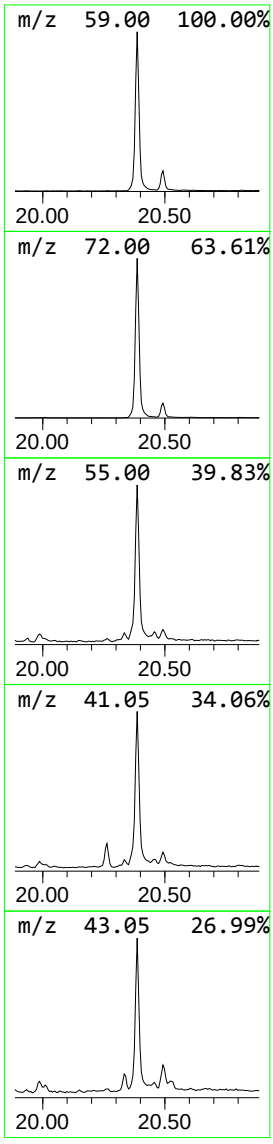
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 9-Octadecenamide, (Z)- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.386	15.42 ng/ul	3428450	Chrysene-d12	21.227

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	98
2	9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	96
3	Tetradecanamide	227	C14H29NO	000638-58-4	72
4	Palmitoleamide	253	C16H31NO	106010-22-4	72
5	Pentanamide, 4-methyl-	115	C6H13NO	001119-29-5	64





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Acq On : 19 Jul 2022 16:41
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Sample : N3753-08
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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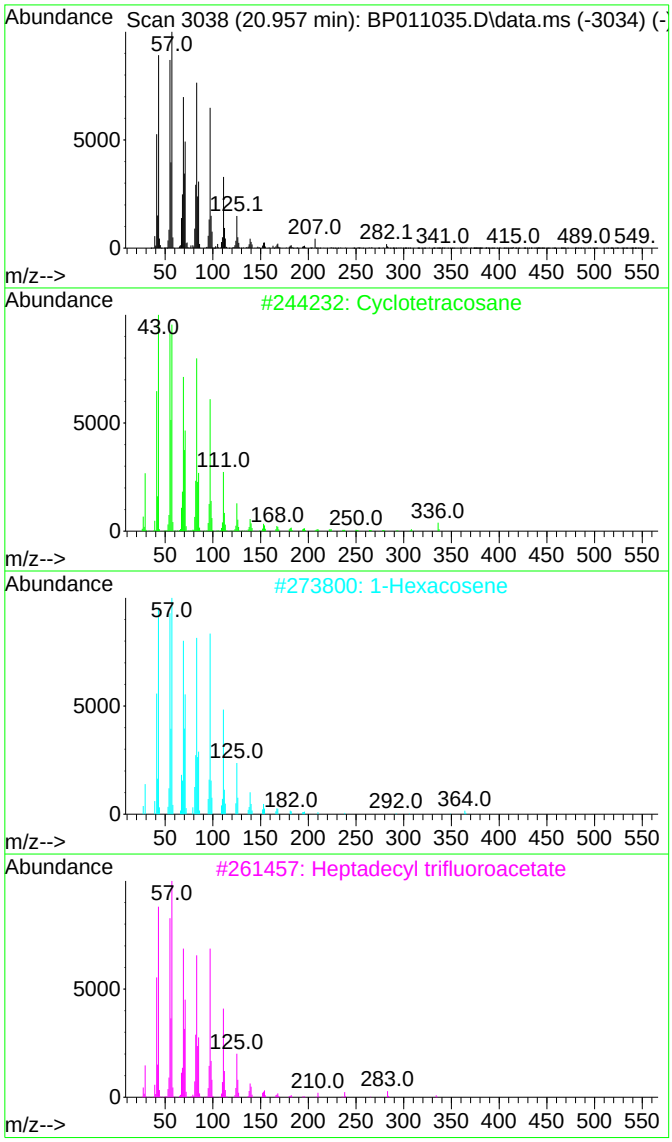
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071422.M
Quant Title : SVOA CALIBRATION

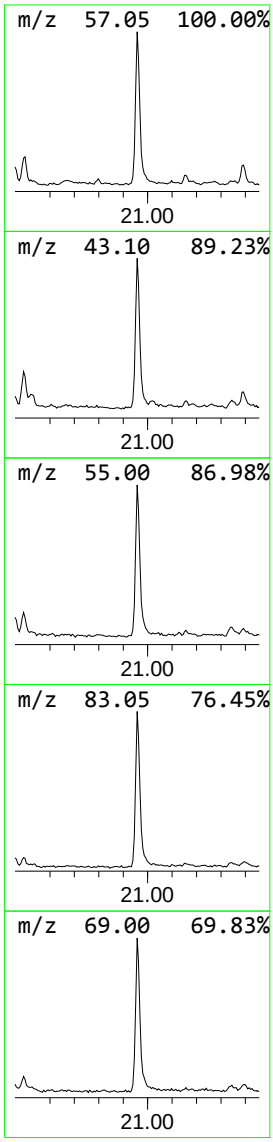
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 11 (DEL) Alkane: Cyclic20.957 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.957	5.91 ng/ul	1314560	Chrysene-d12	21.227

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclotetracosane	336	C24H48	000297-03-0	94
2	1-Hexacosene	364	C26H52	018835-33-1	94
3	Heptadecyl trifluoroacetate	352	C19H35F3O2	1010351-87-0	94
4	Pentafluoropropionic acid, tetra...	360	C17H29F5O2	006222-06-6	94
5	Heptafluorobutyric acid, n-tetra...	410	C18H29F7O2	007365-36-8	91





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Sample : N3753-08
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GBHP9

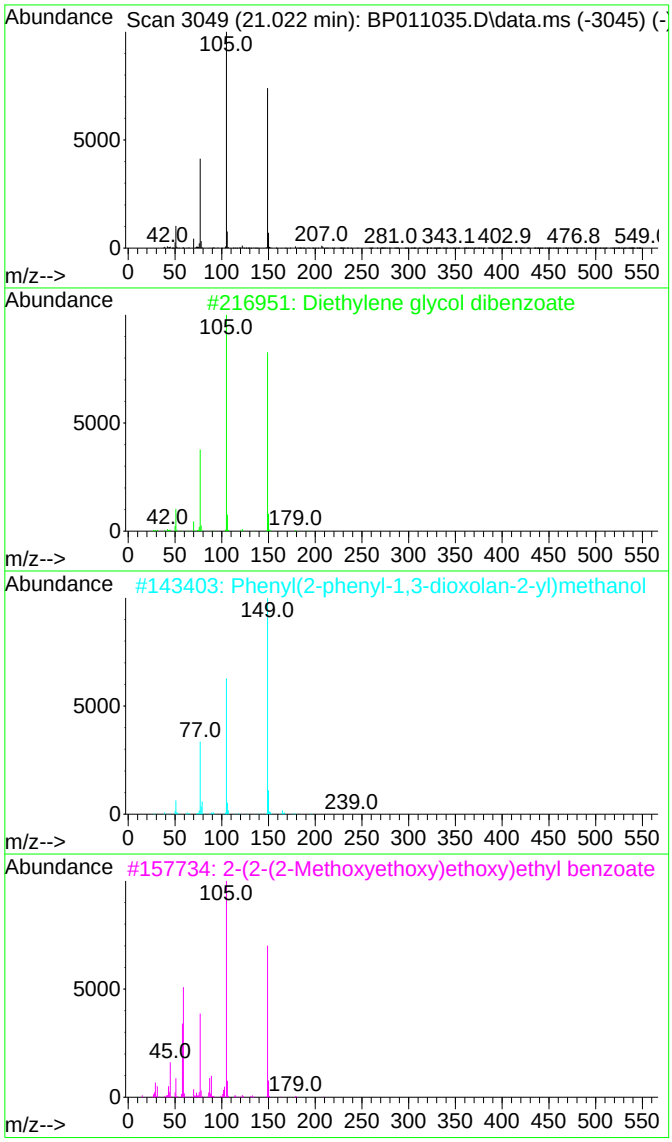
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071422.M
Quant Title : SVOA CALIBRATION

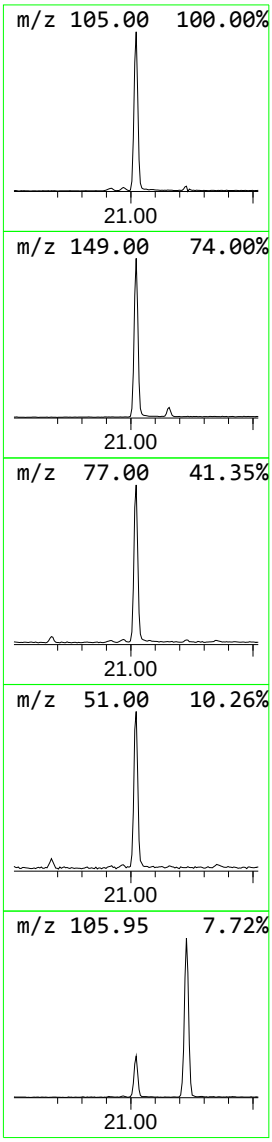
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 12 Diethylene glycol dibenzoate Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.022	5.92 ng/ul	1316190	Chrysene-d12	21.227

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Diethylene glycol dibenzoate	314	C18H18O5	000120-55-8	91
2	Phenyl(2-phenyl-1,3-dioxolan-2-y...	256	C16H16O3	005694-69-9	78
3	2-(2-(2-Methoxyethoxy)ethoxy)eth...	268	C14H20O5	1000366-99-1	74
4	1,3-Dioxolane, 2-(methoxymethyl)...	194	C11H14O3	1000156-71-2	74
5	Diethylene glycol dibenzoate	314	C18H18O5	000120-55-8	72





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Operator : CG/JU
Sample : N3753-08
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071422.M
Quant Title : SVOA CALIBRATION

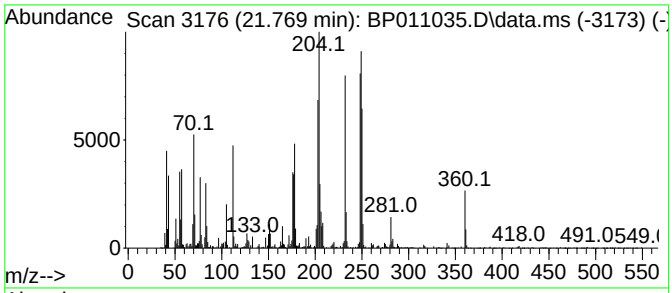
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

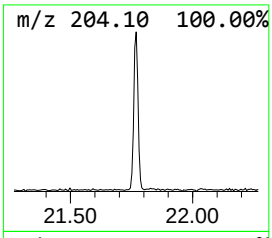
****

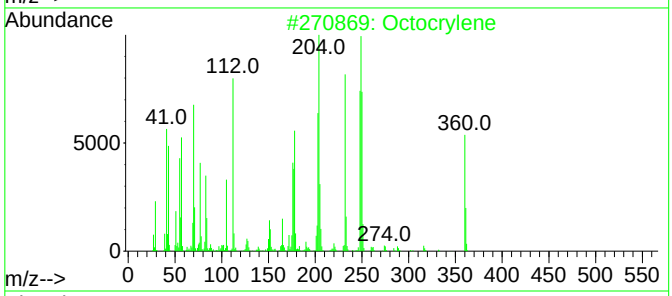
Peak Number 13 Octocrylene Concentration Rank 13

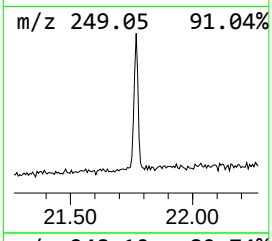
R.T.	EstConc	Area	Relative to ISTD	R.T.
21.769	2.20 ng/ul	490129	Chrysene-d12	21.227

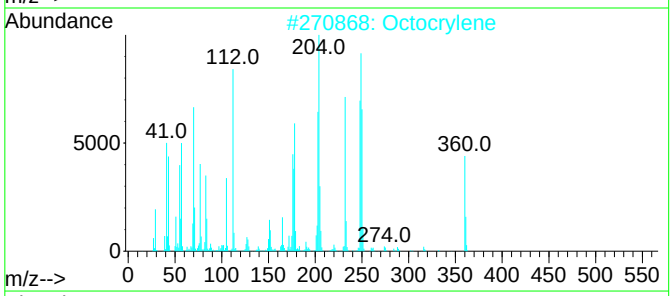
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octocrylene	361	C24H27NO2	006197-30-4	99
2	Octocrylene	361	C24H27NO2	006197-30-4	78
3	Octocrylene	361	C24H27NO2	006197-30-4	50
4	Quinoline-4-carboxamide, 2-pheny...	360	C22H14F2N2O	303133-50-8	30
5	2,5-Dimethyl-7,7,8,8-tetracyano-...	232	C14H8N4	001487-82-7	25

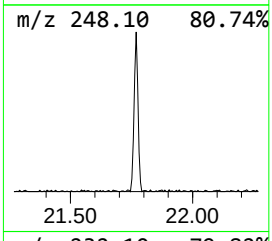


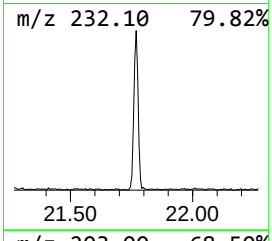


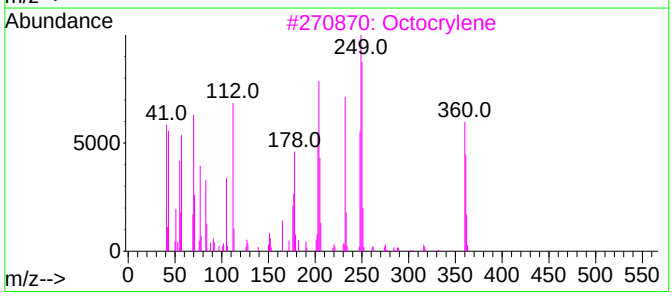


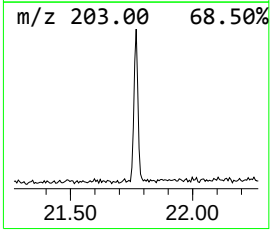












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Data File : BP011035.D
Acq On : 19 Jul 2022 16:41
Operator : CG/JU
Sample : N3753-08
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GBHP9

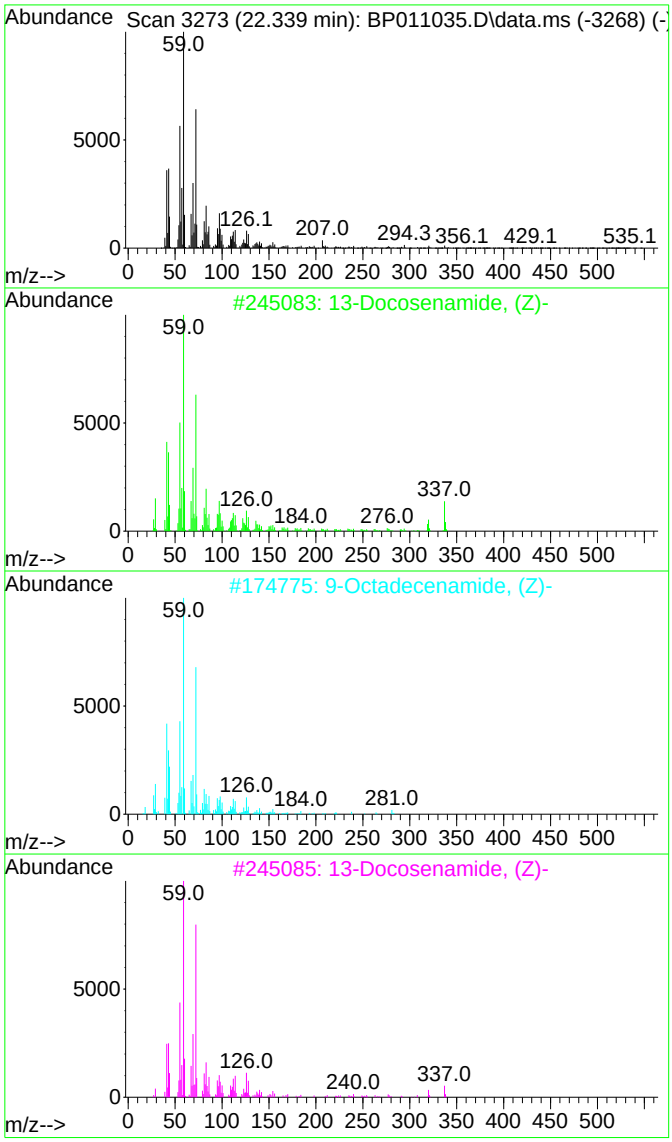
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071422.M
Quant Title : SVOA CALIBRATION

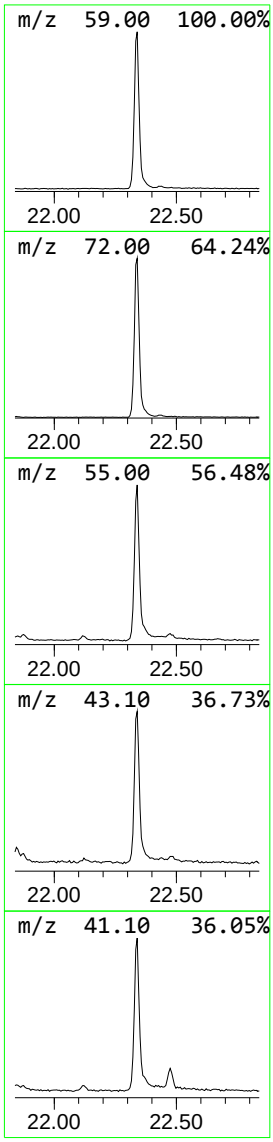
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 14 13-Docosenamide, (Z)- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.339	13.06 ng/ul	2903330	Chrysene-d12	21.227

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	13-Docosenamide, (Z)-	337	C22H43NO	000112-84-5	87
2	9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	80
3	13-Docosenamide, (Z)-	337	C22H43NO	000112-84-5	76
4	9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	74
5	9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	72





Tentatively Identified Compound (LSC) summary

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Data File : BP011035.D
Acq On : 19 Jul 2022 16:41
Operator : CG/JU
Sample : N3753-08
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GBHP9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071422.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
Benzyl alcohol	7.928	2.8	ng/ul	279617	1	7.675	2011650	20.0
Hexanoic acid, ...	8.987	6.2	ng/ul	620824	1	7.675	2011650	20.0
Ethanol, 2-phen...	10.840	33.6	ng/ul	4724190	2	10.469	2812720	20.0
1-Phenoxypropan...	11.228	2.1	ng/ul	290190	2	10.469	2812720	20.0
N-n-Butylphthal...	16.004	26.2	ng/ul	4855470	4	17.104	3708760	20.0
7,9-Di-tert-but...	17.792	3.1	ng/ul	576237	4	17.104	3708760	20.0
(DEL) Alkane: S...	18.863	4.5	ng/ul	842472	4	17.104	3708760	20.0
Bis(2-ethylhexy...	19.169	2.4	ng/ul	528343	5	21.227	4446270	20.0
unknown-01	20.263	2.5	ng/ul	555740	5	21.227	4446270	20.0
9-Octadecenamid...	20.386	15.4	ng/ul	3428450	5	21.227	4446270	20.0
(DEL) Alkane: C...	20.957	5.9	ng/ul	1314560	5	21.227	4446270	20.0
Diethylene glyc...	21.022	5.9	ng/ul	1316190	5	21.227	4446270	20.0
Octocrylene	21.769	2.2	ng/ul	490129	5	21.227	4446270	20.0
13-Docosenamide...	22.339	13.1	ng/ul	2903330	5	21.227	4446270	20.0