LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080824\
Data File : BG062320.D
Acq On : 8 Aug 2024 5:12
Operator : MA/JU
Sample : P3437-12
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
E05X9

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\SFAM-EPA-BG080524.MA.M
Title : SVOA CALIBRATION

Signal : TIC: BG062320.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.420	19	22	34	rVB5	12143	23384	1.28%	0.109%
2	3.684	63	67	73	rBV	10427	16824	0.92%	0.078%
3	3.825	87	91	106	rVV	66765	120761	6.60%	0.562%
4	6.040	463	468	474	rVB	7870	11241	0.61%	0.052%
5	6.892	608	613	619	rBV3	5445	12315	0.67%	0.057%
6	7.109	644	650	660	rVB	74083	133114	7.27%	0.620%
7	7.285	673	680	692	rVB	149573	248624	13.58%	1.158%
8	7.491	708	715	722	rBV	175034	303998	16.60%	1.416%
9	7.550	722	725	730	rVV2	5989	10085	0.55%	0.047%
10	7.955	786	794	802	rVV	188315	328673	17.95%	1.531%
11	8.014	802	804	810	rVB2	5542	8596	0.47%	0.040%
12	8.648	904	912	923	rBV	187251	317417	17.34%	1.479%
13	9.106	982	990	1001	rVV	186863	323406	17.66%	1.506%
14	9.670	1081	1086	1094	rVB2	31553	49903	2.73%	0.232%
15	9.829	1105	1113	1123	rBV	146673	258522	14.12%	1.204%
16	10.364	1196	1204	1214	rBV	282115	494666	27.02%	2.304%
17	10.751	1262	1270	1275	rBV	303453	528931	28.89%	2.464%
18	10.798	1275	1278	1284	rVV	41364	67426	3.68%	0.314%
19	10.875	1284	1291	1305	rVB	220732	404524	22.09%	1.884%
20	11.280	1355	1360	1365	rBV5	11090	17332	0.95%	0.081%
21	11.486	1388	1395	1405	rBV2	84558	154683	8.45%	0.721%
22	12.332	1535	1539	1544	rVB4	4803	8271	0.45%	0.039%
23	12.408	1544	1552	1559	rBV	29659	52974	2.89%	0.247%
24	12.519	1564	1571	1576	rBV	22519	35851	1.96%	0.167%
25	12.625	1583	1589	1594	rVB	21492	34073	1.86%	0.159%
26	13.412	1717	1723	1730	rVB5	7614	13966	0.76%	0.065%
27	13.982	1813	1820	1827	rVB	564740	799559	43.67%	3.724%
28	14.094	1833	1839	1843	rBV2	8605	12722	0.69%	0.059%
29	14.270	1856	1869	1878	rVV	576645	954697	52.14%	4.447%
30	14.576	1914	1921	1927	rBV	569373	864745	47.23%	4.028%
31	14.634	1927	1931	1937	rVB	50176	77318	4.22%	0.360%
32	14.693	1937	1941	1945	rBV	6664	9672	0.53%	0.045%
33	14.752	1945	1951	1956	rBV	72394	133531	7.29%	0.622%
34	14.881	1969	1973	1975	rBV	7106	8931	0.49%	0.042%
35	14.969	1982	1988	1994	rBV	19847	32871	1.80%	0.153%
36	15.316	2037	2047	2053	rVV	529409	733552	40.07%	3.417%

LSC Area Percent Report

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Sample : P3437-12
Misc :
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Instrument :
BNA_G
ClientSampleId :
E05X9

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\SFAM-EPA-BG080524.MA.M
Title : SVOA CALIBRATION

37	15.404	2053	2062	2066	rVV8	9554	28409	1.55%	0.132%
38	15.439	2066	2068	2076	rVB7	8808	13082	0.71%	0.061%
39	15.568	2081	2090	2095	rBV	823558	1259566	68.80%	5.867%
40	15.621	2095	2099	2102	rVV2	27081	39218	2.14%	0.183%
41	15.668	2102	2107	2114	rVV	177774	258197	14.10%	1.203%
42	15.739	2115	2119	2124	rVB7	7175	9839	0.54%	0.046%
43	15.938	2150	2153	2156	rVB6	6214	8224	0.45%	0.038%
44	16.120	2180	2184	2189	rVB3	10693	18534	1.01%	0.086%
45	16.244	2196	2205	2206	rBV5	10499	19919	1.09%	0.093%
46	16.285	2208	2212	2217	rVB5	13334	28309	1.55%	0.132%
47	16.467	2239	2243	2246	rBV5	5989	10885	0.59%	0.051%
48	16.720	2280	2286	2290	rBV4	12281	19820	1.08%	0.092%
49	16.990	2328	2332	2337	rVB7	6180	11248	0.61%	0.052%
50	17.090	2345	2349	2353	rVV4	17840	23998	1.31%	0.112%
51	17.131	2353	2356	2361	rVB6	7422	13630	0.74%	0.063%
52	17.248	2372	2376	2380	rBV5	13929	17745	0.97%	0.083%
53	17.313	2380	2387	2392	rBV	733099	1067402	58.30%	4.972%
54	17.354	2392	2394	2398	rVV2	28030	44529	2.43%	0.207%
55	17.413	2398	2404	2413	rVV	945944	1396215	76.26%	6.503%
56	17.501	2414	2419	2424	rVB3	16287	29760	1.63%	0.139%
57	17.707	2445	2454	2460	rBV	40836	75214	4.11%	0.350%
58	17.859	2477	2480	2484	rVB6	4830	7221	0.39%	0.034%
59	17.942	2490	2494	2500	rBV7	9736	17084	0.93%	0.080%
60	17.994	2500	2503	2509	rVB7	6836	10881	0.59%	0.051%
61	18.218	2534	2541	2559	rBV	329860	551209	30.11%	2.567%
62	18.881	2650	2654	2657	rBV5	9730	12747	0.70%	0.059%
63	19.022	2673	2678	2681	rBV7	10497	18262	1.00%	0.085%
64	19.075	2682	2687	2696	rVV6	30817	68130	3.72%	0.317%
65	19.146	2696	2699	2704	rVB4	21185	28115	1.54%	0.131%
66	19.257	2716	2718	2724	rVB4	16699	25156	1.37%	0.117%
67	19.334	2727	2731	2732	rBV	14431	17280	0.94%	0.080%
68	19.487	2753	2757	2767	rBV2	98840	144198	7.88%	0.672%
69	19.639	2779	2783	2786	rBV5	9778	12048	0.66%	0.056%
70	19.692	2786	2792	2801	rBV	1166090	1734307	94.73%	8.078%
71	20.021	2844	2848	2849	rBV3	7959	9699	0.53%	0.045%
72	20.144	2866	2869	2877	rVB2	14914	24318	1.33%	0.113%
73	20.215	2878	2881	2884	rVV3	18277	20827	1.14%	0.097%
74	20.256	2886	2888	2891	rVB3	14491	14301	0.78%	0.067%
75	21.231	3049	3054	3061	rBV	261353	386376	21.10%	1.800%
76	21.472	3093	3095	3100	rVB6	10111	11686	0.64%	0.054%

LSC Area Percent Report

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Misc :
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Integration Parameters: LSCINT.P
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Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG080524.MA.M
Title : SVOA CALIBRATION

77	21.554	3102	3109	3125	rVB	746785	1223030	66.80%	5.697%
78	22.236	3221	3225	3232	rVB7	23408	40551	2.21%	0.189%
79	22.306	3232	3237	3238	rBV4	11656	18312	1.00%	0.085%
80	22.347	3239	3244	3246	rVV2	52920	89959	4.91%	0.419%
81	22.377	3246	3249	3255	rVB2	56884	94738	5.17%	0.441%
82	22.524	3268	3274	3282	rVB4	49188	93414	5.10%	0.435%
83	22.659	3291	3297	3302	rBV2	99807	201667	11.01%	0.939%
84	22.717	3304	3307	3317	rVB4	70541	130427	7.12%	0.608%
85	23.041	3360	3362	3370	rVB5	15202	25185	1.38%	0.117%
86	23.229	3388	3394	3401	rBV2	134924	247489	13.52%	1.153%
87	23.828	3490	3496	3501	rBV2	35349	71757	3.92%	0.334%
88	24.427	3587	3598	3616	rVB	677616	1830858	100.00%	8.528%
89	24.638	3623	3634	3646	rBV	478373	1298104	70.90%	6.046%
90	24.744	3648	3652	3662	rVB5	25883	56006	3.06%	0.261%
91	25.837	3831	3838	3844	rBV3	31350	82078	4.48%	0.382%
92	25.943	3848	3856	3868	rVB4	38477	127344	6.96%	0.593%
93	26.783	3990	3999	4010	rVB4	40585	124951	6.82%	0.582%
94	26.953	4024	4028	4029	rBV4	6738	8710	0.48%	0.041%
95	27.147	4055	4061	4070	rVB4	20025	59117	3.23%	0.275%
96	27.881	4179	4186	4198	rVB4	18803	69186	3.78%	0.322%
97	28.721	4319	4329	4339	rVB9	26802	108929	5.95%	0.507%
98	28.927	4356	4364	4366	rBV6	26557	59931	3.27%	0.279%
99	29.761	4497	4506	4515	rBV6	23953	90679	4.95%	0.422%
100	30.031	4543	4552	4566	rVB6	47333	201626	11.01%	0.939%

Sum of corrected areas: 21468794

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080824\
Data File : BG062320.D
Acq On : 8 Aug 2024 5:12
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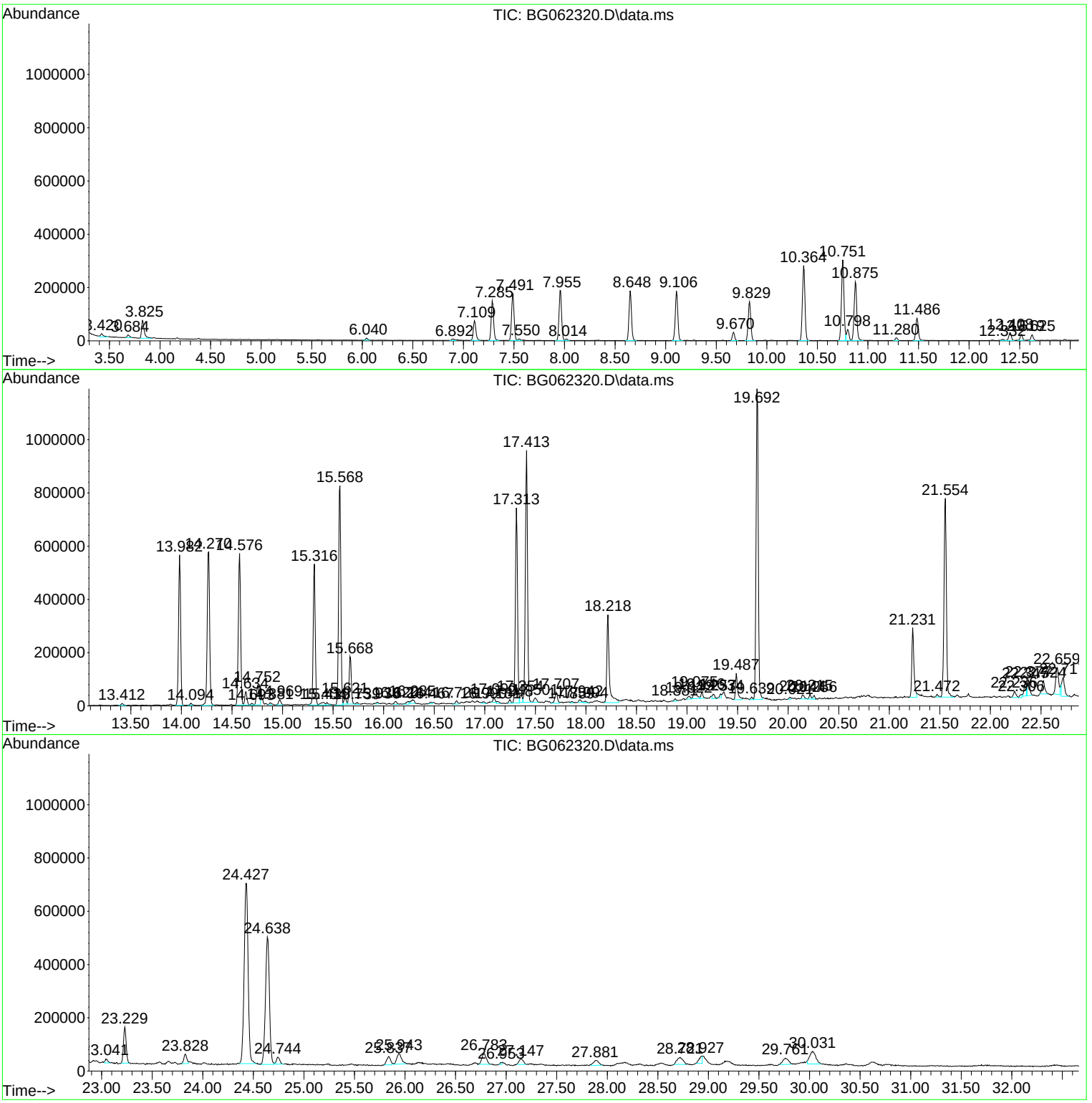
Instrument :
BNA_G
ClientSampleId :
E05X9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG080524.MA.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080824\
Data File : BG062320.D
Acq On : 8 Aug 2024 5:12
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG080524.MA.M
Quant Title : SVOA CALIBRATION

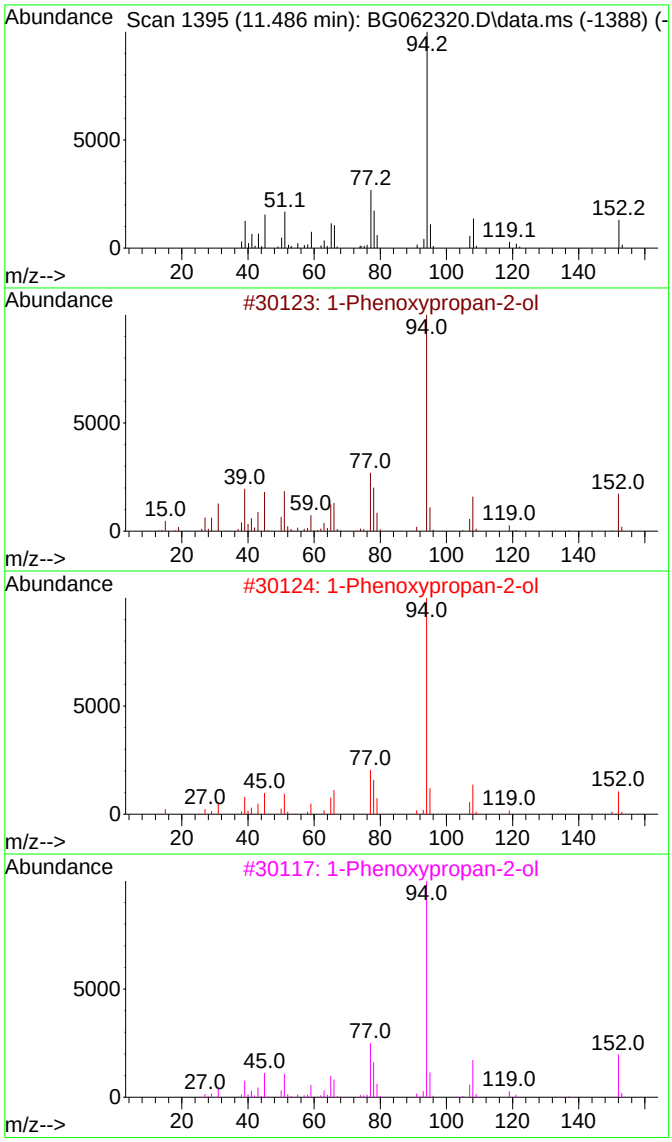
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

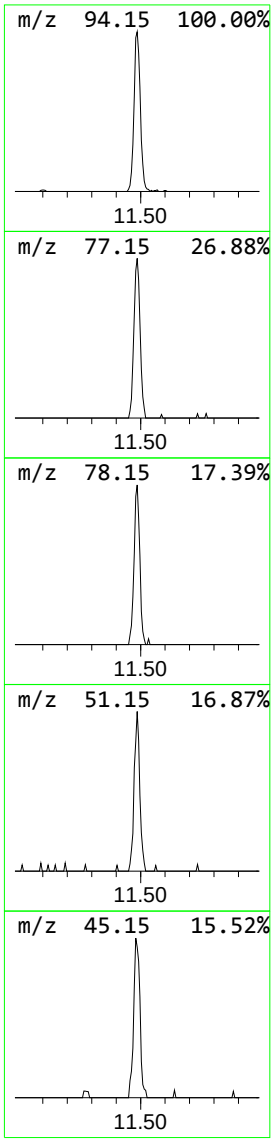
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.486	5.85 ng/ul	154683	Naphthalene-d8	10.751

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	96
3	1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	95
4	1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	87
5	1-Propanol, 3-phenoxy-	152	C9H12O2	006180-61-6	80





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080824\
Data File : BG062320.D
Acq On : 8 Aug 2024 5:12
Operator : MA/JU
Sample : P3437-12
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
BNA_G
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG080524.MA.M
Quant Title : SVOA CALIBRATION

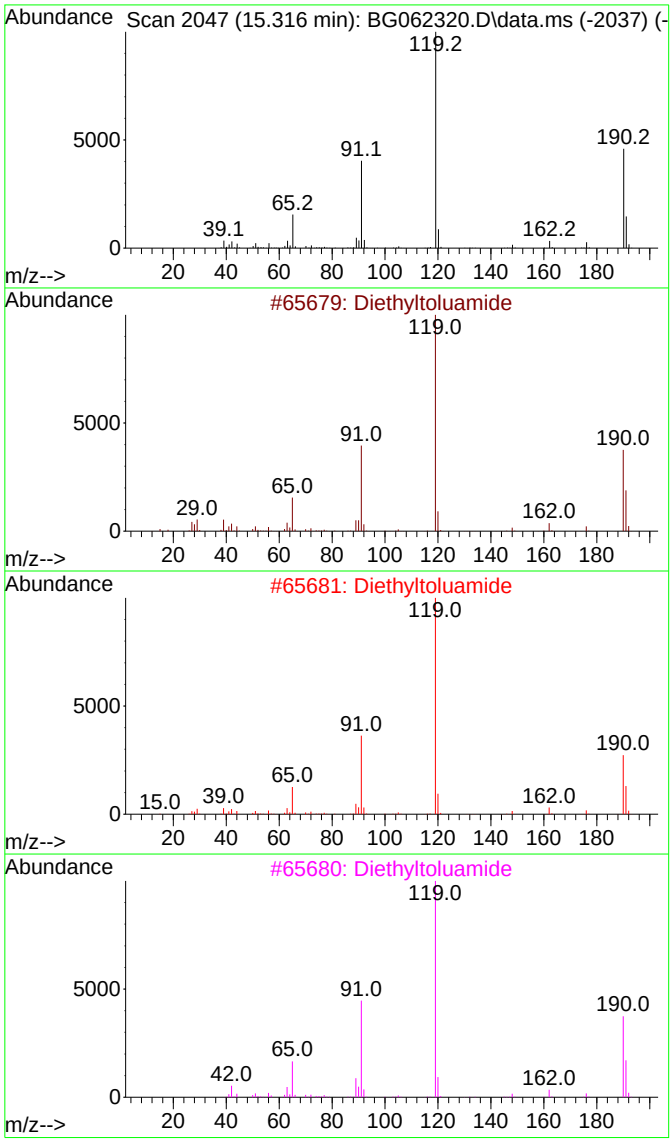
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

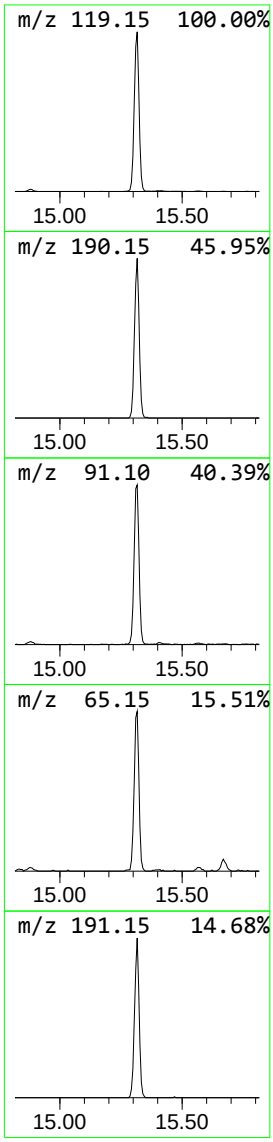
****

Peak Number 2 Diethyltoluamide Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.316	16.97 ng/ul	733552	Acenaphthene-d10	14.576

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Diethyltoluamide	191	C12H17NO	000134-62-3	96
2	Diethyltoluamide	191	C12H17NO	000134-62-3	95
3	Diethyltoluamide	191	C12H17NO	000134-62-3	94
4	Benzamide, N,N-diethyl-4-methyl-	191	C12H17NO	002728-05-4	93
5	Benzamide, 3-methyl-N-methyl-N-p...	191	C12H17NO	1000421-46-2	81





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080824\
Data File : BG062320.D
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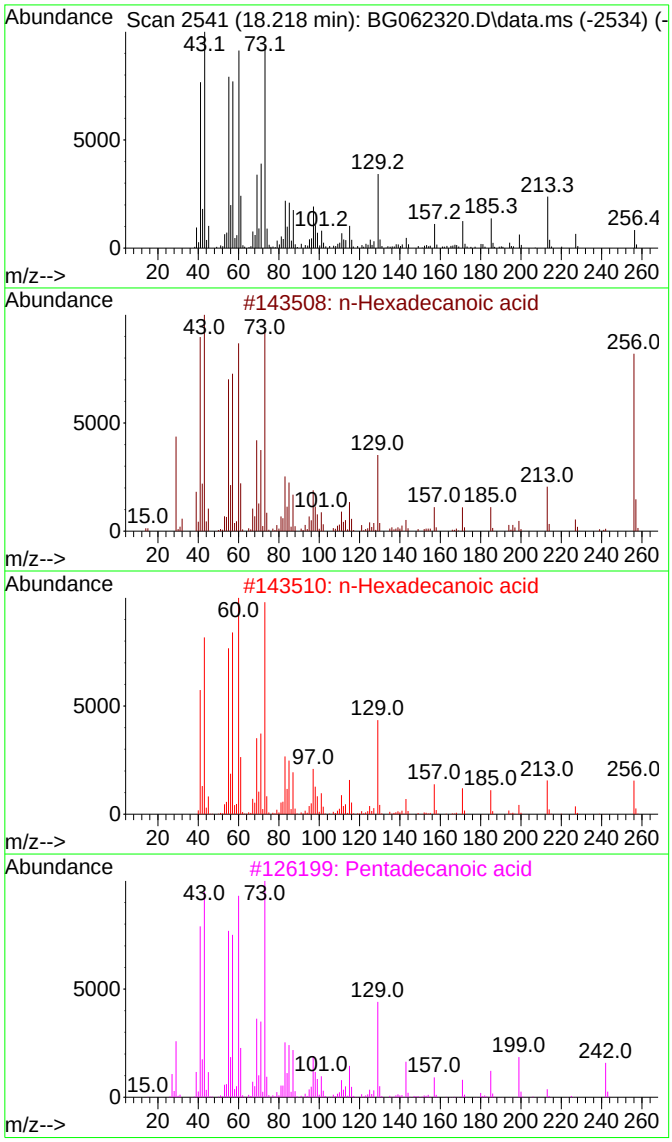
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG080524.MA.M
Quant Title : SVOA CALIBRATION

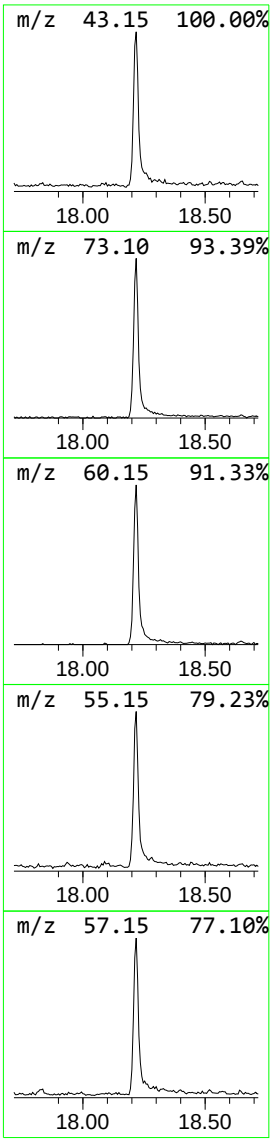
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 n-Hexadecanoic acid Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.218	10.33 ng/ul	551209	Phenanthrene-d10	17.313

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
2	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	91
3	Pentadecanoic acid	242	C15H30O2	001002-84-2	90
4	Tridecanoic acid	214	C13H26O2	000638-53-9	89
5	Tetradecanoic acid	228	C14H28O2	000544-63-8	87





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Sample : P3437-12
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG080524.MA.M
Quant Title : SVOA CALIBRATION

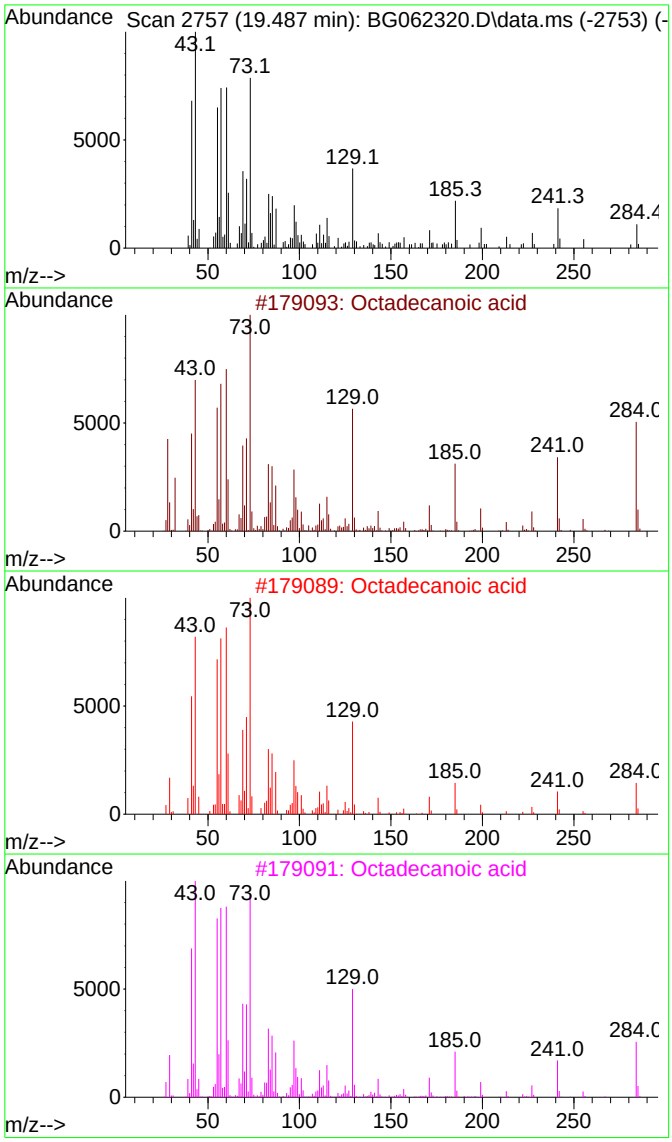
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

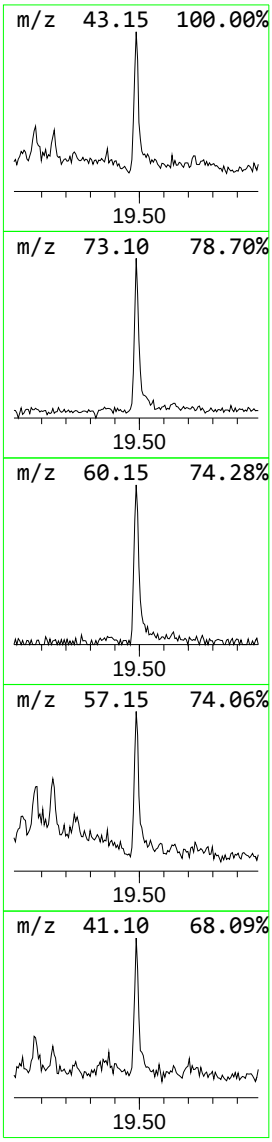
****

Peak Number 4 Octadecanoic acid Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.486	2.36 ng/ul	144198	Chrysene-d12	21.554

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecanoic acid	284	C18H36O2	000057-11-4	98
2	Octadecanoic acid	284	C18H36O2	000057-11-4	95
3	Octadecanoic acid	284	C18H36O2	000057-11-4	95
4	Octadecanoic acid	284	C18H36O2	000057-11-4	87
5	Octadecanoic acid	284	C18H36O2	000057-11-4	83





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Data File : BG062320.D
Acq On : 8 Aug 2024 5:12
Operator : MA/JU
Sample : P3437-12
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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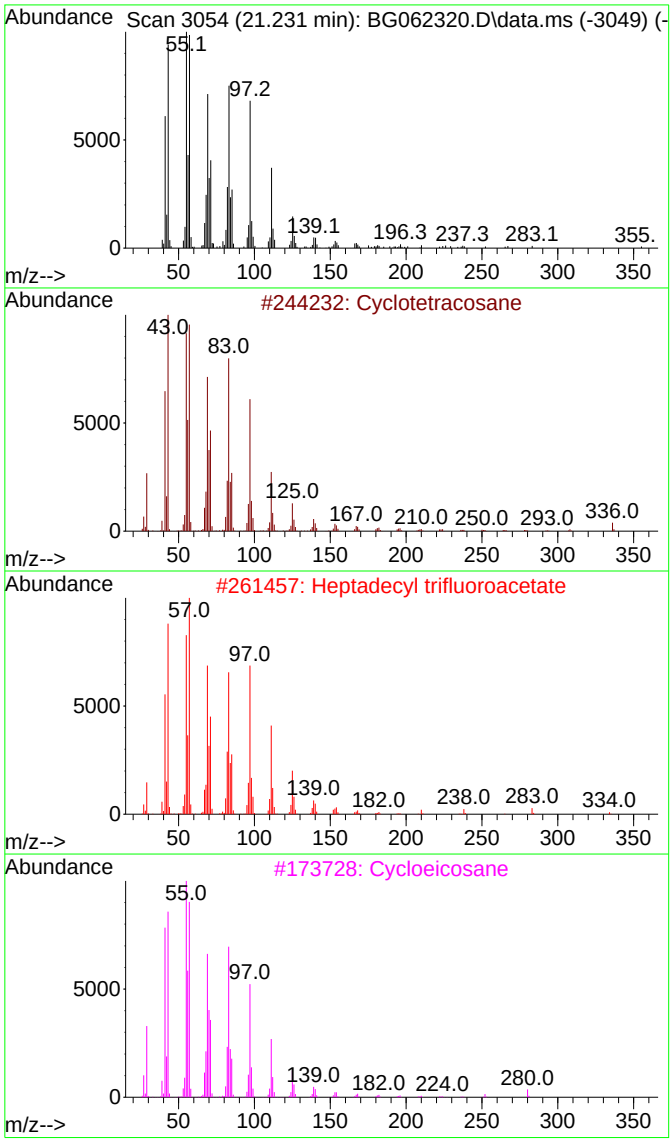
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG080524.MA.M
Quant Title : SVOA CALIBRATION

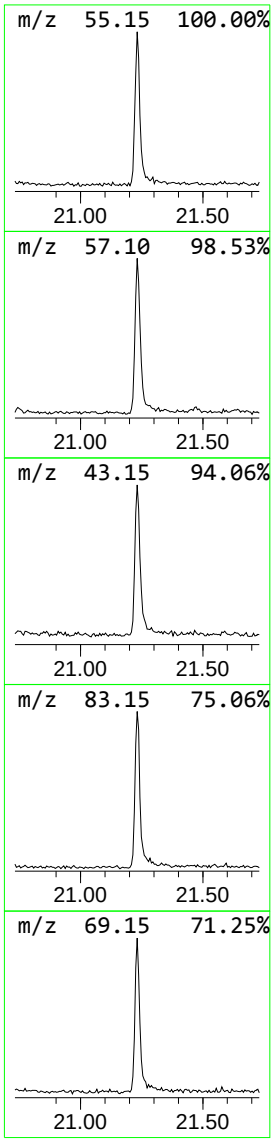
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 (DEL) Alkane: Cyclic21.231 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.231	6.32 ng/ul	386376	Chrysene-d12	21.554

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclotetracosane	336	C24H48	000297-03-0	95
2	Heptadecyl trifluoroacetate	352	C19H35F3O2	1010351-87-0	94
3	Cycloeicosane	280	C20H40	000296-56-0	91
4	n-Tetracosanol-1	354	C24H50O	000506-51-4	91
5	Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	91





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Data File : BG062320.D
Acq On : 8 Aug 2024 5:12
Operator : MA/JU
Sample : P3437-12
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
E05X9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG080524.MA.M
Quant Title : SVOA CALIBRATION

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

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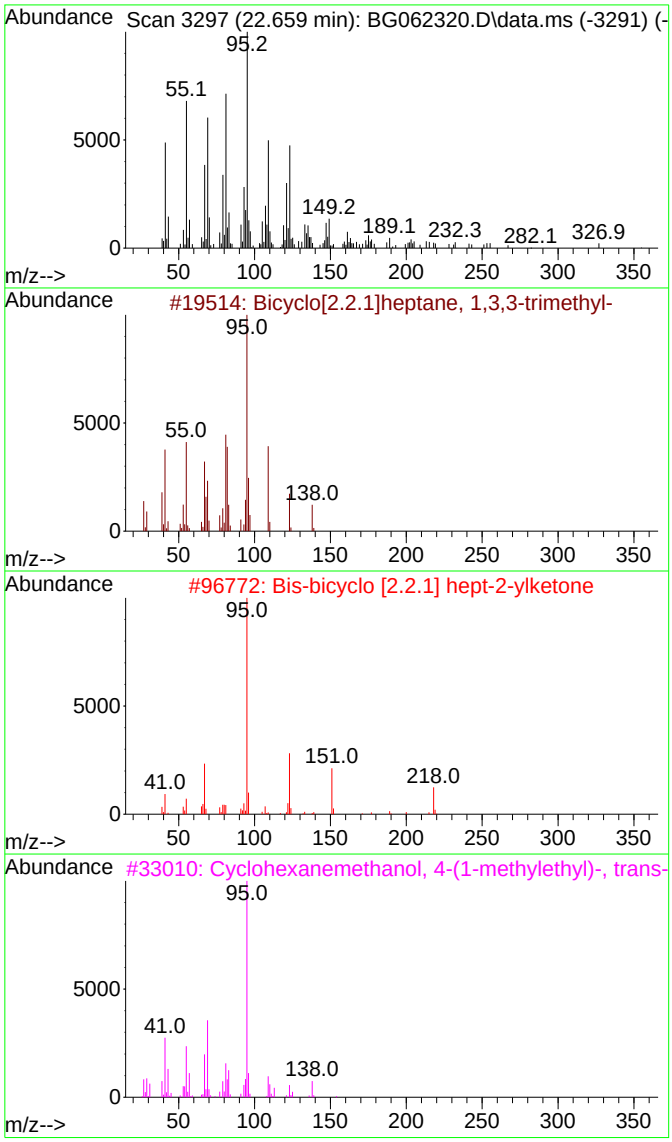
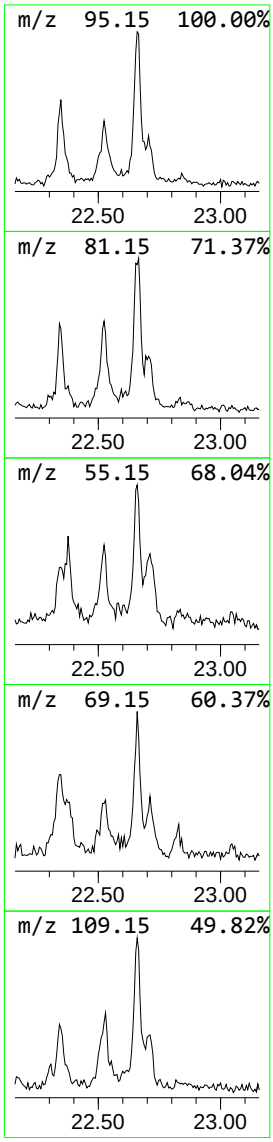
Peak Number 6 unknown-01 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.659	3.30 ng/ul	201667	Chrysene-d12	21.554

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Bicyclo[2.2.1]heptane, 1,3,3-tri...	138	C10H18	006248-88-0	47
2	Bis-bicyclo [2.2.1] hept-2-ylketone	218	C15H22O	1010222-21-8	43
3	Cyclohexanemethanol, 4-(1-methyl...	156	C10H20O	013674-19-6	38
4	Silane, diethyldifluoro-	124	C4H10F2Si	000358-06-5	38
5	5,7-Octadien-4-one, 2,6-dimethyl...	152	C10H16O	003588-18-9	35





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080824\
Data File : BG062320.D
Acq On : 8 Aug 2024 5:12
Operator : MA/JU
Sample : P3437-12
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
E05X9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG080524.MA.M
Quant Title : SVOA CALIBRATION

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

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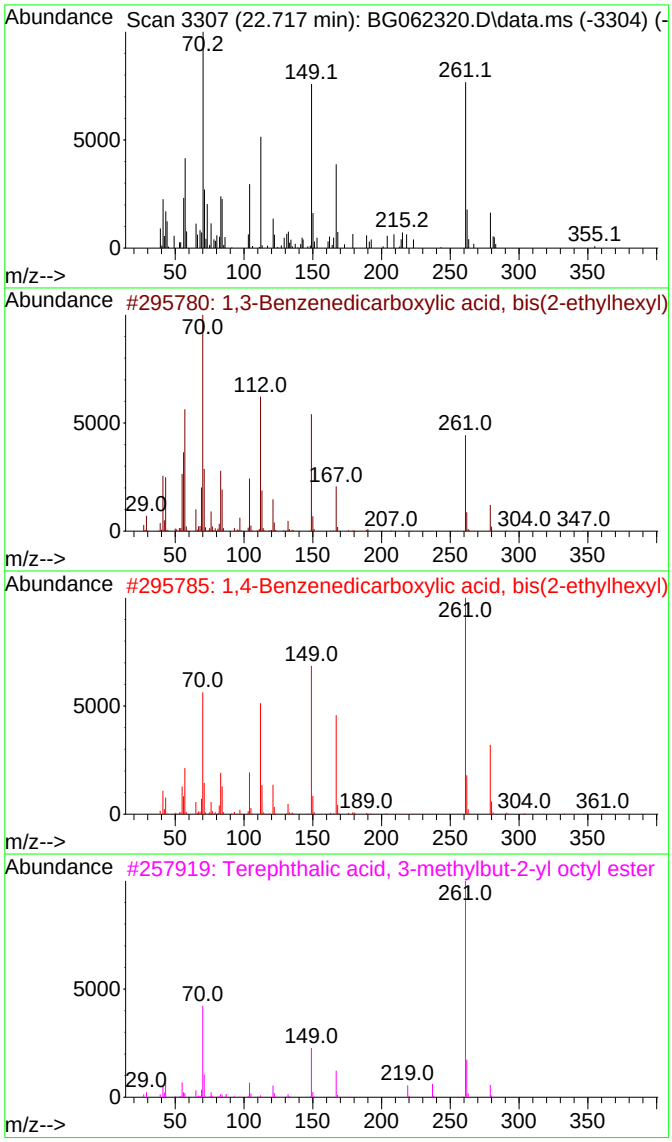
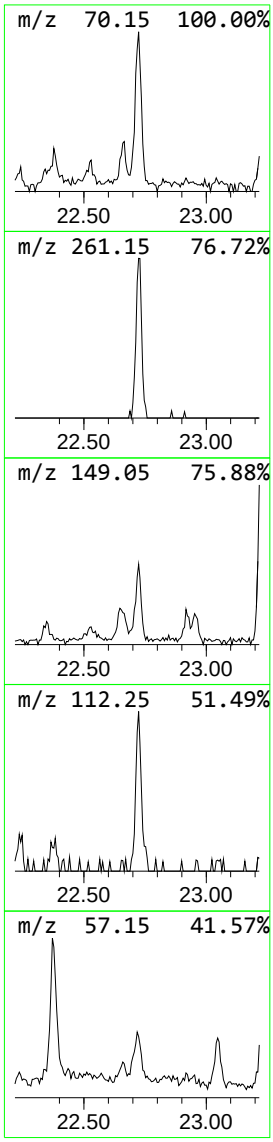
Peak Number 7 1,3-Benzenedicarboxylic aci... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.717	2.13 ng/ul	130427	Chrysene-d12	21.554

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,3-Benzenedicarboxylic acid, bi...	390	C24H38O4	000137-89-3	64
2	1,4-Benzenedicarboxylic acid, bi...	390	C24H38O4	006422-86-2	47
3	Terephthalic acid, 3-methylbut-2...	348	C21H32O4	1000356-27-6	43
4	1,3-Benzenedicarboxylic acid, bi...	390	C24H38O4	000137-89-3	38
5	Terephthalic acid, 2-bromo-4-flu...	450	C22H24BrFO4	1000323-71-2	35





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Data File : BG062320.D
Acq On : 8 Aug 2024 5:12
Operator : MA/JU
Sample : P3437-12
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
E05X9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG080524.MA.M
Quant Title : SVOA CALIBRATION

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

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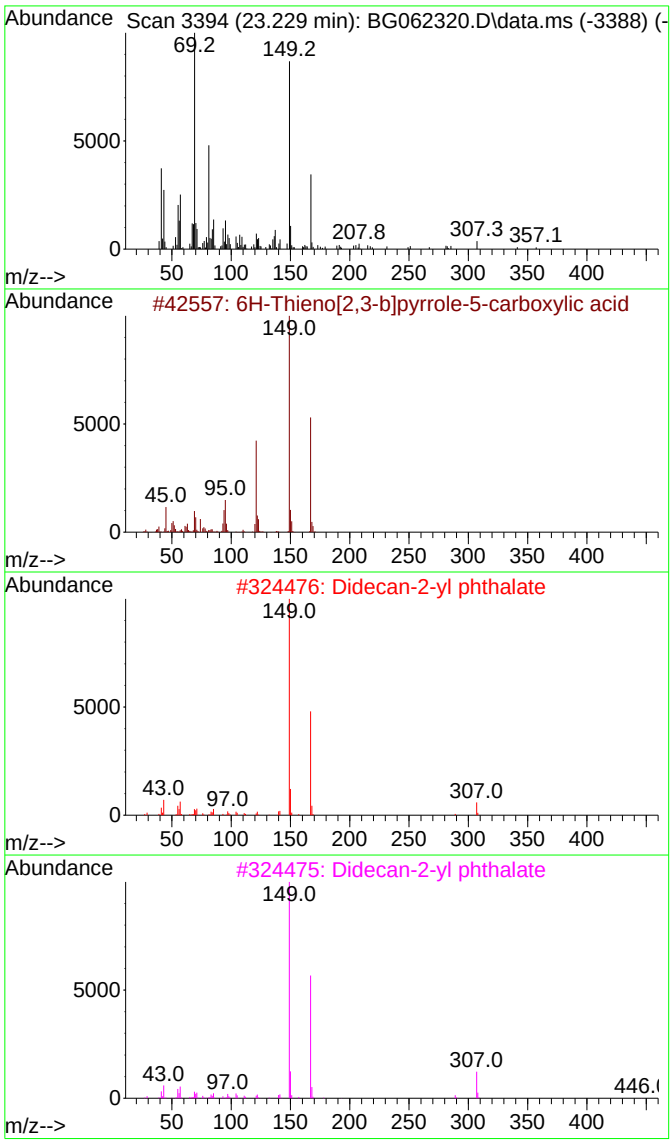
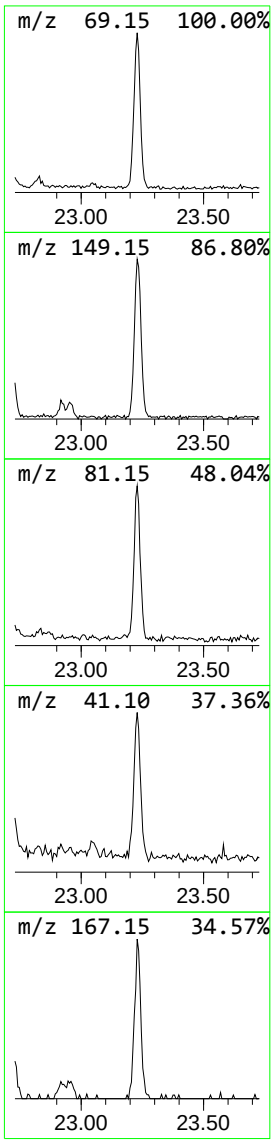
Peak Number 8 unknown-02 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.229	3.81 ng/ul	247489	Perylene-d12	24.638

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	6H-Thieno[2,3-b]pyrrole-5-carbox...	167	C7H5NO2S	051856-25-8	43
2	Didecan-2-yl phthalate	446	C28H46O4	028029-89-2	38
3	Didecan-2-yl phthalate	446	C28H46O4	028029-89-2	38
4	Bis(2-ethylhexyl) phthalate	390	C24H38O4	000117-81-7	35
5	Phthalic acid, 5-methylhex-2-yl ...	502	C32H54O4	1000371-07-7	35





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080824\
Data File : BG062320.D
Acq On : 8 Aug 2024 5:12
Operator : MA/JU
Sample : P3437-12
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
E05X9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG080524.MA.M
Quant Title : SVOA CALIBRATION

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

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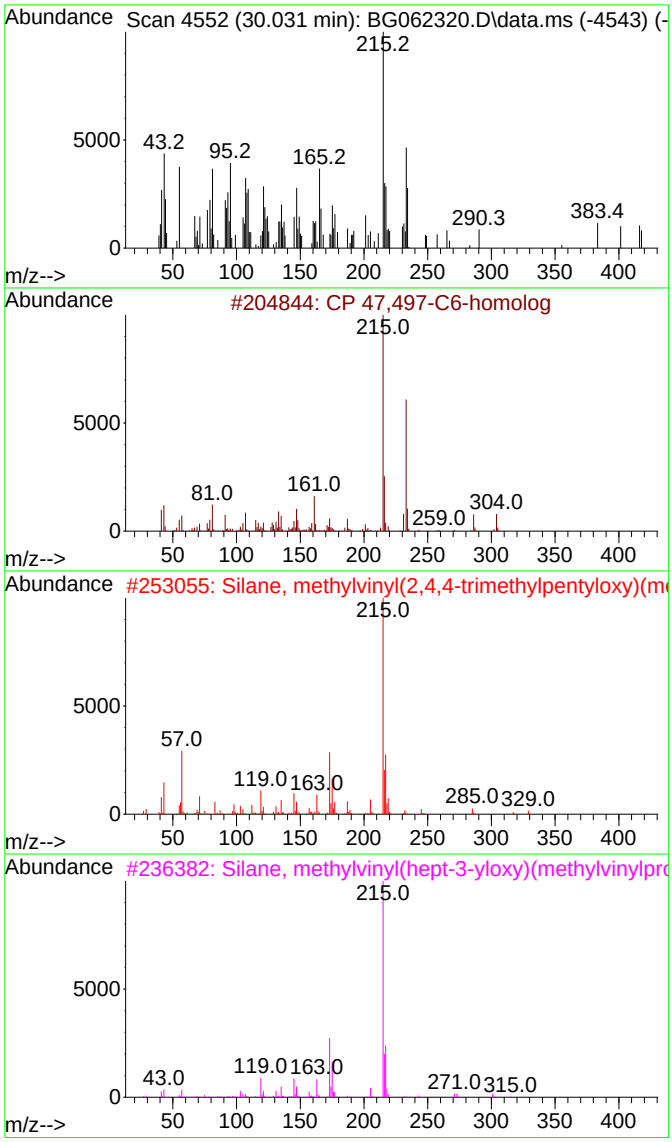
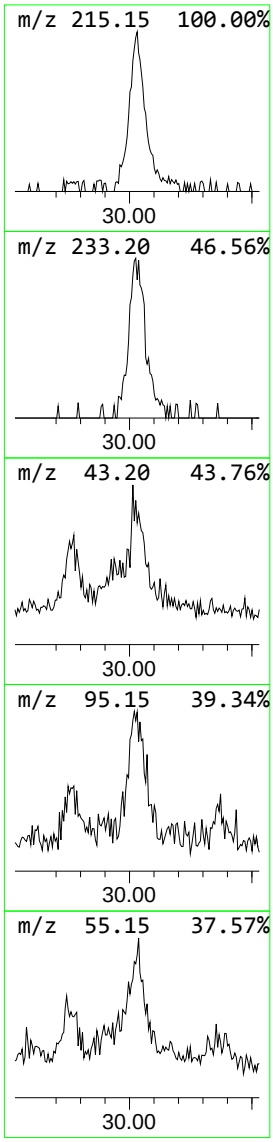
Peak Number 9 unknown-03 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
30.031	3.11 ng/ul	201626	Perylene-d12	24.638

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	CP 47,497-C6-homolog	304	C20H32O2	070435-06-2	37
2	Silane, methylvinyl(2,4,4-trimet...	344	C17H36O3Si2	1000416-58-8	27
3	Silane, methylvinyl(hept-3-yloxy...	330	C16H34O3Si2	1010421-69-1	27
4	Silane, methylvinyl(hex-2-yloxy)...	316	C15H32O3Si2	1000421-59-6	27
5	5-(2-Methyloctan-2-yl)-3-(1R,3R-...	332	C22H36O2	070434-92-3	25





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080824\
Data File : BG062320.D
Acq On : 8 Aug 2024 5:12
Operator : MA/JU
Sample : P3437-12
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
E05X9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG080524.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
1-Phenoxypropan...	11.486	5.8	ng/ul	154683	2	10.751	528931	20.0
Diethyltoluamide	15.316	17.0	ng/ul	733552	3	14.576	864745	20.0
n-Hexadecanoic ...	18.218	10.3	ng/ul	551209	4	17.313	1067400	20.0
Octadecanoic acid	19.486	2.4	ng/ul	144198	5	21.554	1223030	20.0
(DEL) Alkane: C...	21.231	6.3	ng/ul	386376	5	21.554	1223030	20.0
unknown-01	22.659	3.3	ng/ul	201667	5	21.554	1223030	20.0
1,3-Benzenedica...	22.717	2.1	ng/ul	130427	5	21.554	1223030	20.0
unknown-02	23.229	3.8	ng/ul	247489	6	24.638	1298100	20.0
unknown-03	30.031	3.1	ng/ul	201626	6	24.638	1298100	20.0