LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062224\
Data File : BG061660.D
Acq On : 22 Jun 2024 22:36
Operator : MA/JU
Sample : P2776-21
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
A4C74

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-BG062024.MA.M
Title : SVOA CALIBRATION

Signal : TIC: BG061660.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.470	26	31	36	rVB2	20356	33174	1.55%	0.114%
2	3.740	72	77	83	rBV	54377	85714	4.01%	0.293%
3	3.811	88	89	94	rBV4	4299	6681	0.31%	0.023%
4	3.887	99	102	113	rBV	277041	457374	21.41%	1.565%
5	4.193	151	154	157	rBV5	4029	5360	0.25%	0.018%
6	4.234	157	161	167	rVB5	5600	10123	0.47%	0.035%
7	4.686	236	238	241	rVB3	5148	4969	0.23%	0.017%
8	4.792	252	256	261	rVB4	4374	7834	0.37%	0.027%
9	5.150	313	317	320	rBV3	9249	11416	0.53%	0.039%
10	5.679	405	407	411	rBV4	4977	4530	0.21%	0.016%
11	5.720	411	414	416	rBV3	4523	4978	0.23%	0.017%
12	6.132	479	484	493	rVV3	18869	33819	1.58%	0.116%
13	6.337	514	519	522	rVB3	3198	4970	0.23%	0.017%
14	6.596	560	563	567	rBV2	4191	6251	0.29%	0.021%
15	6.754	583	590	596	rVV	296939	480886	22.51%	1.646%
16	7.060	634	642	648	rVB3	27022	43979	2.06%	0.150%
17	7.219	662	669	682	rVB	397652	680779	31.87%	2.330%
18	7.401	692	700	709	rBV	255712	433184	20.28%	1.482%
19	7.512	712	719	728	rBV	740858	1242503	58.17%	4.252%
20	7.601	728	734	741	rVB	346650	589646	27.61%	2.018%
21	7.659	741	744	747	rBV2	10611	13142	0.62%	0.045%
22	7.859	775	778	780	rBV2	3720	4501	0.21%	0.015%
23	7.988	795	800	803	rVB4	5178	8113	0.38%	0.028%
24	8.076	807	815	821	rVV	304582	526744	24.66%	1.802%
25	8.135	821	825	828	rVB3	11488	19246	0.90%	0.066%
26	8.300	846	853	857	rBV5	21855	38293	1.79%	0.131%
27	8.358	857	863	870	rVB4	20103	33165	1.55%	0.113%
28	8.770	925	933	940	rBV	470380	787347	36.86%	2.694%
29	9.240	1004	1013	1020	rBV	368525	645332	30.21%	2.208%
30	9.293	1020	1022	1026	rVV5	5383	7866	0.37%	0.027%
31	9.398	1035	1040	1046	rVB4	6066	10516	0.49%	0.036%
32	9.510	1055	1059	1060	rBV3	3760	5018	0.23%	0.017%
33	9.798	1100	1108	1113	rBV3	42581	75252	3.52%	0.258%
34	9.962	1129	1136	1145	rVB2	305273	549323	25.72%	1.880%
35	10.227	1179	1181	1185	rVB3	6159	6658	0.31%	0.023%
36	10.497	1219	1227	1238	rBV	486893	873447	40.89%	2.989%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062224\
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Acq On : 22 Jun 2024 22:36
Operator : MA/JU
Sample : P2776-21
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
A4C74

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-BG062024.MA.M
Title : SVOA CALIBRATION

37	10.656	1247	1254	1259	rBV2	27488	47868	2.24%	0.164%
38	10.891	1286	1294	1301	rVV	491818	883591	41.37%	3.024%
39	11.020	1305	1316	1328	rBV	424096	790299	37.00%	2.704%
40	11.114	1330	1332	1337	rVV3	3664	4873	0.23%	0.017%
41	11.279	1357	1360	1364	rBV4	5112	6987	0.33%	0.024%
42	11.414	1379	1383	1388	rBV5	14686	25213	1.18%	0.086%
43	11.625	1413	1419	1426	rVB2	77940	140038	6.56%	0.479%
44	11.854	1456	1458	1461	rVB4	4544	4569	0.21%	0.016%
45	11.895	1461	1465	1468	rBV3	3549	5979	0.28%	0.020%
46	11.948	1472	1474	1478	rBV2	3034	4446	0.21%	0.015%
47	12.166	1508	1511	1515	rVB4	4172	4469	0.21%	0.015%
48	12.295	1531	1533	1537	rVB4	5218	6261	0.29%	0.021%
49	12.460	1557	1561	1568	rVB3	10531	23021	1.08%	0.079%
50	12.512	1568	1570	1573	rBV2	3406	4687	0.22%	0.016%
51	12.759	1610	1612	1616	rBV2	4618	4644	0.22%	0.016%
52	12.912	1637	1638	1642	rBV2	3930	6168	0.29%	0.021%
53	13.076	1662	1666	1669	rVB5	3573	5076	0.24%	0.017%
54	13.523	1739	1742	1744	rBV3	7765	7580	0.35%	0.026%
55	13.817	1787	1792	1798	rVB4	19651	29696	1.39%	0.102%
56	14.028	1823	1828	1833	rBV5	15720	26464	1.24%	0.091%
57	14.111	1835	1842	1850	rVV	769462	1157134	54.17%	3.960%
58	14.210	1858	1859	1861	rBV	6079	5431	0.25%	0.019%
59	14.340	1877	1881	1884	rBV6	10018	14120	0.66%	0.048%
60	14.398	1884	1891	1898	rVB	936509	1454340	68.09%	4.977%
61	14.539	1912	1915	1920	rVB4	19613	26977	1.26%	0.092%
62	14.645	1931	1933	1934	rBV2	6180	4906	0.23%	0.017%
63	14.704	1936	1943	1950	rVV	755029	1215807	56.92%	4.160%
64	14.769	1950	1954	1960	rVB3	40132	60809	2.85%	0.208%
65	14.874	1965	1972	1978	rBV	443727	752829	35.25%	2.576%
66	15.033	1995	1999	2005	rVV6	11050	20618	0.97%	0.071%
67	15.098	2006	2010	2019	rVV9	10915	22331	1.05%	0.076%
68	15.303	2042	2045	2050	rBV5	5294	8770	0.41%	0.030%
69	15.491	2074	2077	2079	rBV4	4189	5804	0.27%	0.020%
70	15.544	2083	2086	2089	rBV4	6676	7017	0.33%	0.024%
71	15.650	2101	2104	2105	rBV2	4086	4777	0.22%	0.016%
72	15.691	2105	2111	2118	rVV	1153284	1725051	80.76%	5.903%
73	15.797	2123	2129	2137	rVV	297587	446482	20.90%	1.528%
74	15.932	2147	2152	2159	rBV10	9857	20366	0.95%	0.070%
75	16.044	2169	2171	2173	rVB3	6808	4347	0.20%	0.015%
76	16.173	2188	2193	2199	rVB3	49409	77969	3.65%	0.267%

LSC Area Percent Report

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Data File : BG061660.D
Acq On : 22 Jun 2024 22:36
Operator : MA/JU
Sample : P2776-21
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
A4C74

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 >
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Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG062024.MA.M
Title : SVOA CALIBRATION

77	16.825	2303	2304	2309	rVB4	11946	14793	0.69%	0.051%
78	17.166	2361	2362	2365	rBV3	6639	5761	0.27%	0.020%
79	17.324	2387	2389	2392	rBV3	6251	6488	0.30%	0.022%
80	17.442	2403	2409	2415	rBV	942643	1432917	67.08%	4.903%
81	17.542	2420	2426	2434	rVV2	1173451	1753674	82.10%	6.001%
82	17.818	2471	2473	2478	rVB6	9256	13390	0.63%	0.046%
83	18.200	2536	2538	2542	rVB6	12083	12341	0.58%	0.042%
84	18.329	2554	2560	2570	rBV2	247109	417711	19.56%	1.429%
85	18.511	2587	2591	2595	rBV7	14021	20502	0.96%	0.070%
86	18.658	2614	2616	2620	rBV4	12950	16803	0.79%	0.057%
87	18.817	2640	2643	2648	rVB	91848	119248	5.58%	0.408%
88	18.958	2665	2667	2671	rBV5	11734	16311	0.76%	0.056%
89	19.263	2715	2719	2723	rBV2	36129	54741	2.56%	0.187%
90	19.487	2749	2757	2763	rBV3	110808	218768	10.24%	0.749%
91	19.598	2773	2776	2784	rBV8	71053	119836	5.61%	0.410%
92	19.821	2808	2814	2825	rBV	1352917	2135978	100.00%	7.309%
93	20.297	2892	2895	2898	rBV4	82794	107349	5.03%	0.367%
94	20.368	2904	2907	2912	rVB6	137054	168452	7.89%	0.576%
95	21.196	3044	3048	3055	rVB2	248163	335412	15.70%	1.148%
96	21.361	3073	3076	3085	rVB	279469	410403	19.21%	1.404%
97	21.637	3120	3123	3128	rVB4	211508	273948	12.83%	0.937%
98	21.708	3130	3135	3141	rVB2	895258	1381891	64.70%	4.729%
99	24.716	3640	3647	3659	rVB3	715933	1886999	88.34%	6.457%
100	24.939	3678	3685	3696	rVB2	543827	1479718	69.28%	5.063%

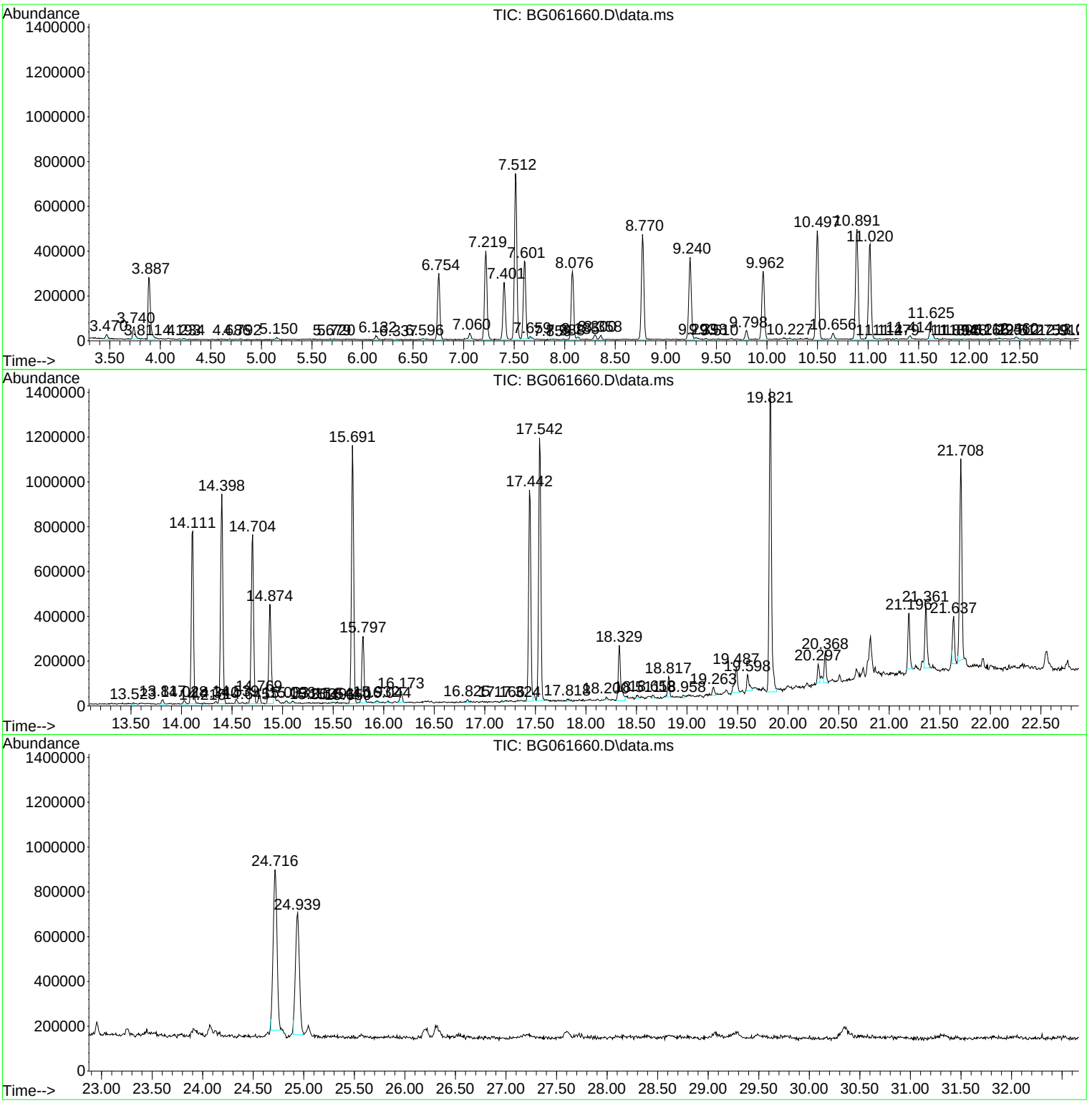
Sum of corrected areas: 29223381

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062224\
Data File : BG061660.D
Acq On : 22 Jun 2024 22:36
Operator : MA/JU
Sample : P2776-21
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
A4C74

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG062024.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\BG062224\
Data File : BG061660.D
Acq On : 22 Jun 2024 22:36
Operator : MA/JU
Sample : P2776-21
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ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG062024.MA.M
Quant Title : SVOA CALIBRATION

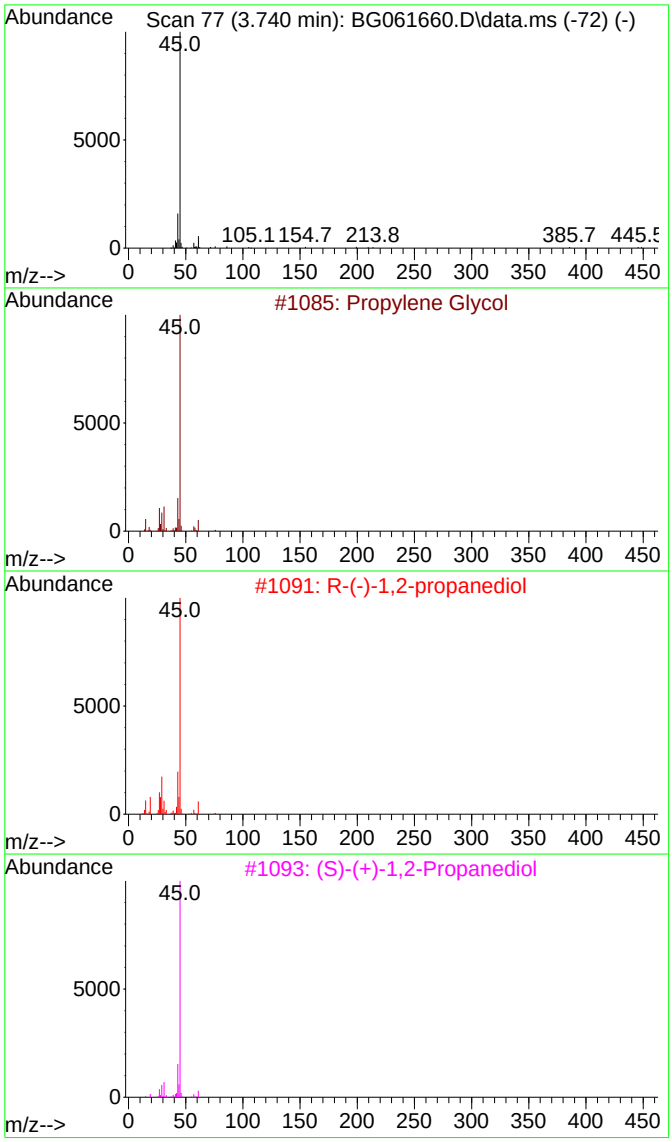
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

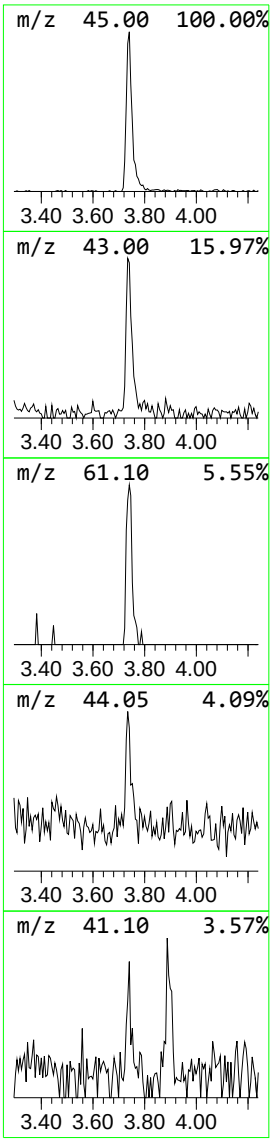
****

Peak Number 1 Propylene Glycol Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.740	3.25 ng/ul	85714	1,4-Dichlorobenzene-d4	8.076

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Propylene Glycol	76	C3H8O2	000057-55-6	56
2	R-(-)-1,2-propanediol	76	C3H8O2	004254-14-2	45
3	(S)-(+)-1,2-Propanediol	76	C3H8O2	004254-15-3	39
4	(S)-(+)-1,2-Propanediol	76	C3H8O2	004254-15-3	10
5	Propylene Glycol	76	C3H8O2	000057-55-6	9





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062224\
Data File : BG061660.D
Acq On : 22 Jun 2024 22:36
Operator : MA/JU
Sample : P2776-21
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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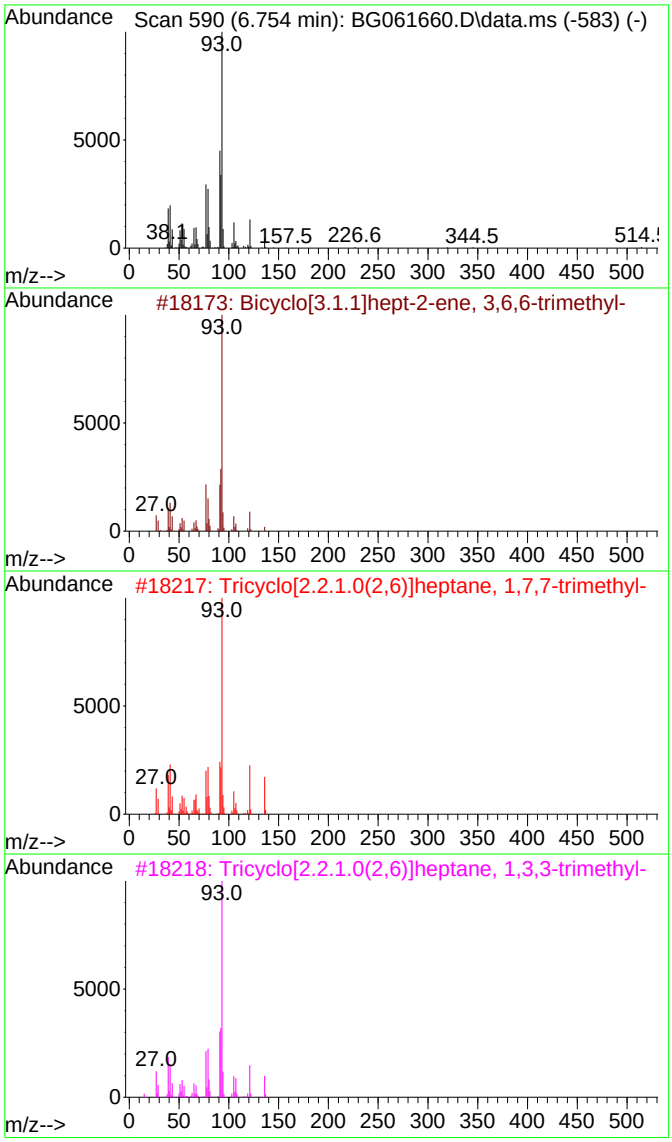
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG062024.MA.M
Quant Title : SVOA CALIBRATION

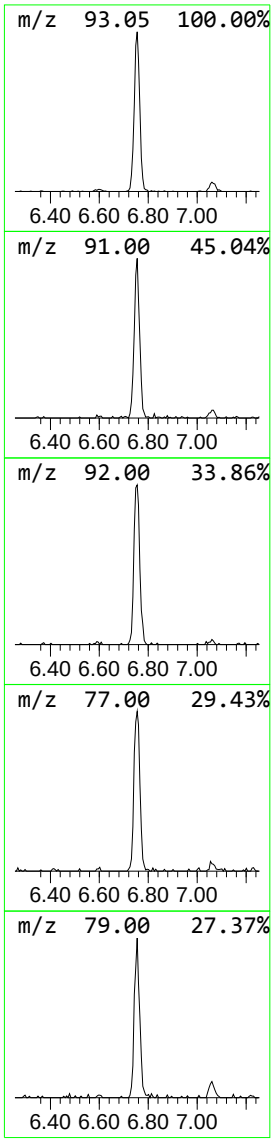
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 Bicyclo[3.1.1]hept-2-ene, 3... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.754	18.26 ng/ul	480886	1,4-Dichlorobenzene-d4	8.076

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Bicyclo[3.1.1]hept-2-ene, 3,6,6-...	136	C10H16	004889-83-2	91
2	Tricyclo[2.2.1.0(2,6)]heptane, 1...	136	C10H16	000508-32-7	91
3	Tricyclo[2.2.1.0(2,6)]heptane, 1...	136	C10H16	000488-97-1	91
4	Tricyclo[2.2.1.0(2,6)]heptane, 1...	136	C10H16	000508-32-7	91
5	(1R)-2,6,6-Trimethylbicyclo[3.1....	136	C10H16	007785-70-8	90





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062224\
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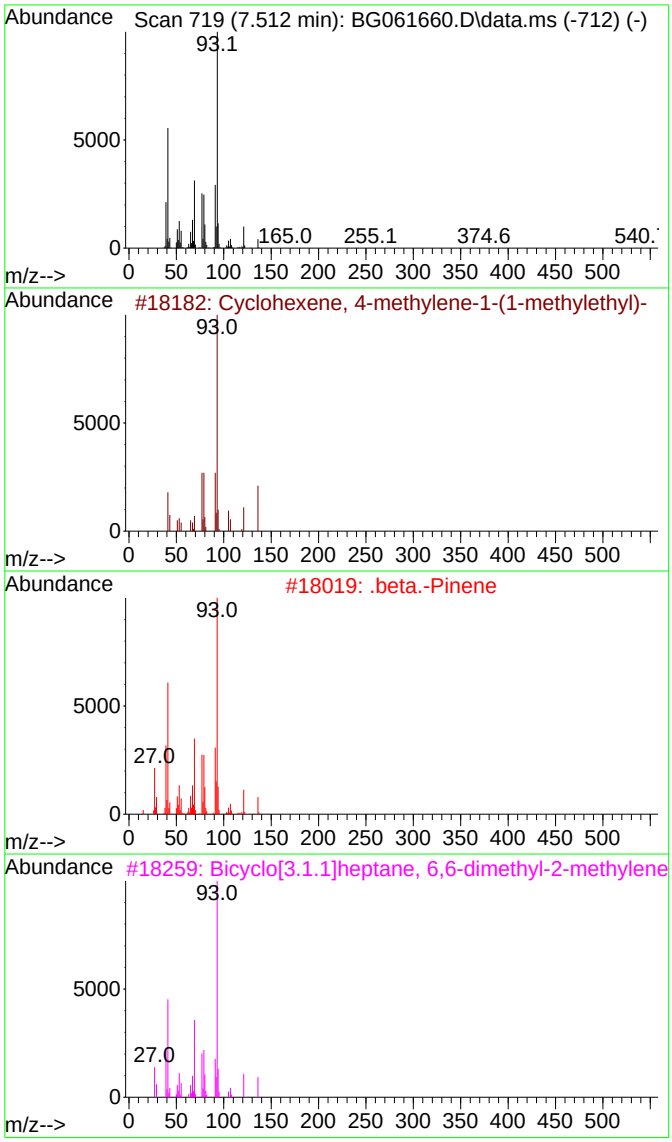
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG062024.MA.M
Quant Title : SVOA CALIBRATION

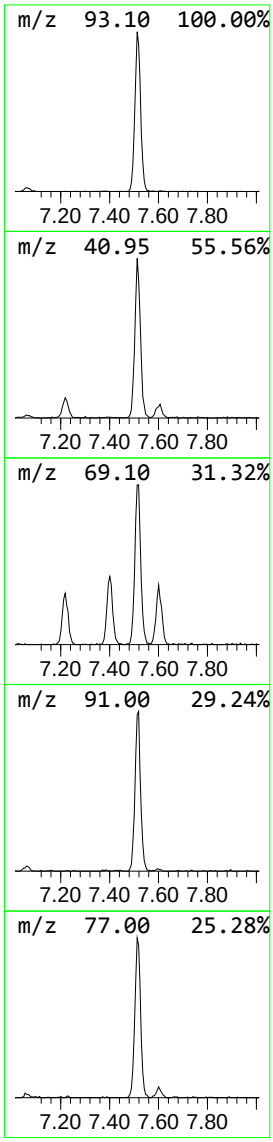
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Cyclohexene, 4-methylene-1-... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.512	47.18 ng/ul	1242500	1,4-Dichlorobenzene-d4	8.076

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexene, 4-methylene-1-(1-me...	136	C10H16	000099-84-3	91
2	.beta.-Pinene	136	C10H16	000127-91-3	91
3	Bicyclo[3.1.1]heptane, 6,6-dimet...	136	C10H16	018172-67-3	91
4	(1R)-2,6,6-Trimethylbicyclo[3.1....	136	C10H16	007785-70-8	87
5	.beta.-Pinene	136	C10H16	000127-91-3	83





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062224\
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Acq On : 22 Jun 2024 22:36
Operator : MA/JU
Sample : P2776-21
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG062024.MA.M
Quant Title : SVOA CALIBRATION

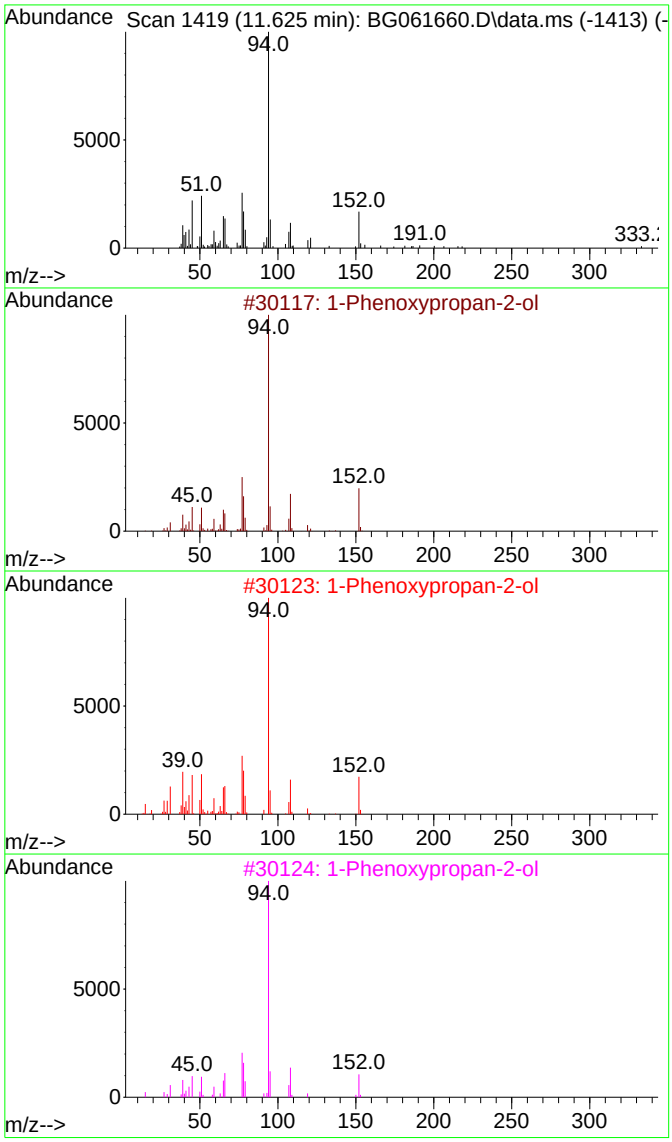
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

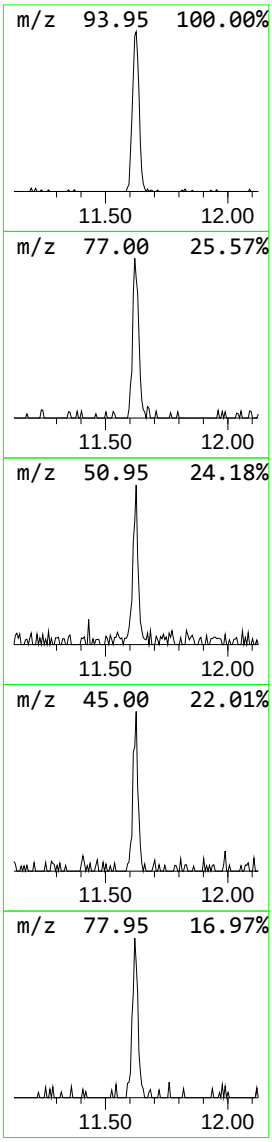
****

Peak Number 4 1-Phenoxypropan-2-ol Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.625	3.17 ng/ul	140038	Naphthalene-d8	10.891

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	93
2	1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	93
3	1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	90
4	1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	81
5	1-Propanol, 3-phenoxy-	152	C9H12O2	006180-61-6	64





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062224\
Data File : BG061660.D
Acq On : 22 Jun 2024 22:36
Operator : MA/JU
Sample : P2776-21
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
A4C74

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG062024.MA.M
Quant Title : SVOA CALIBRATION

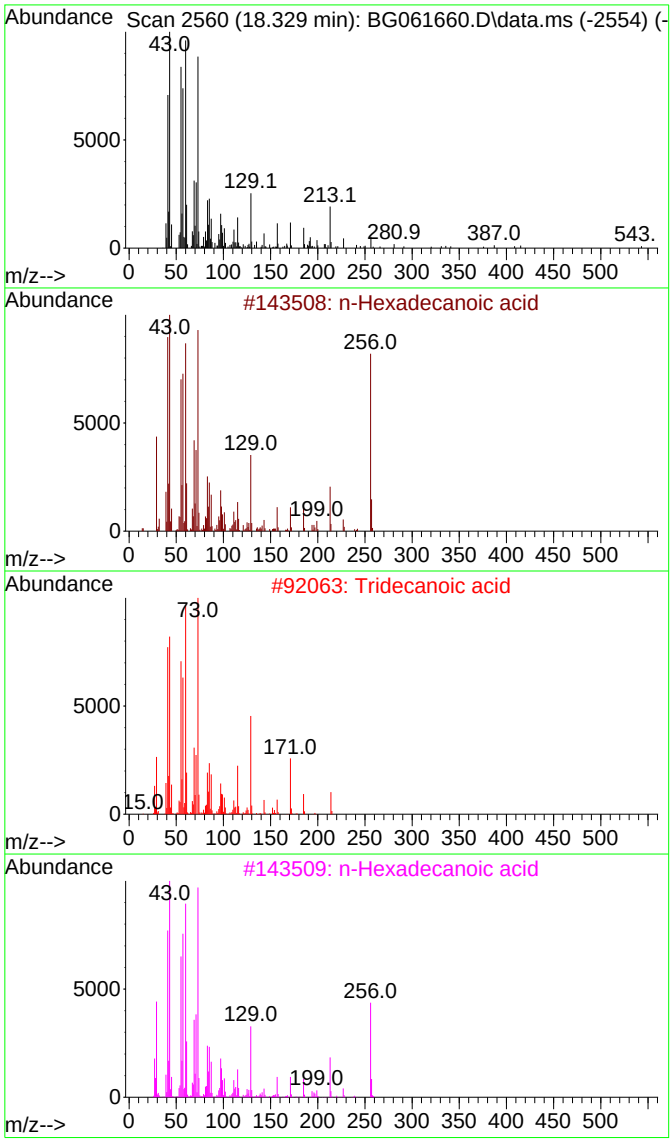
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

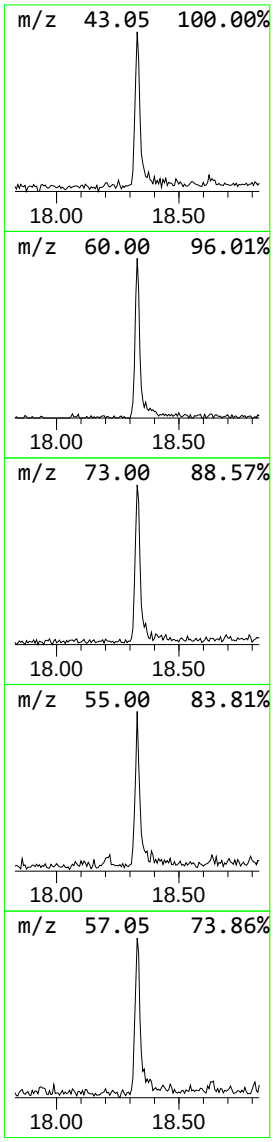
****

Peak Number 5 n-Hexadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.329	5.83 ng/ul	417711	Phenanthrene-d10	17.442

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
2	Tridecanoic acid	214	C13H26O2	000638-53-9	93
3	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	93
4	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	83
5	Tridecanoic acid	214	C13H26O2	000638-53-9	83





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Data File : BG061660.D
Acq On : 22 Jun 2024 22:36
Operator : MA/JU
Sample : P2776-21
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
A4C74

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG062024.MA.M
Quant Title : SVOA CALIBRATION

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

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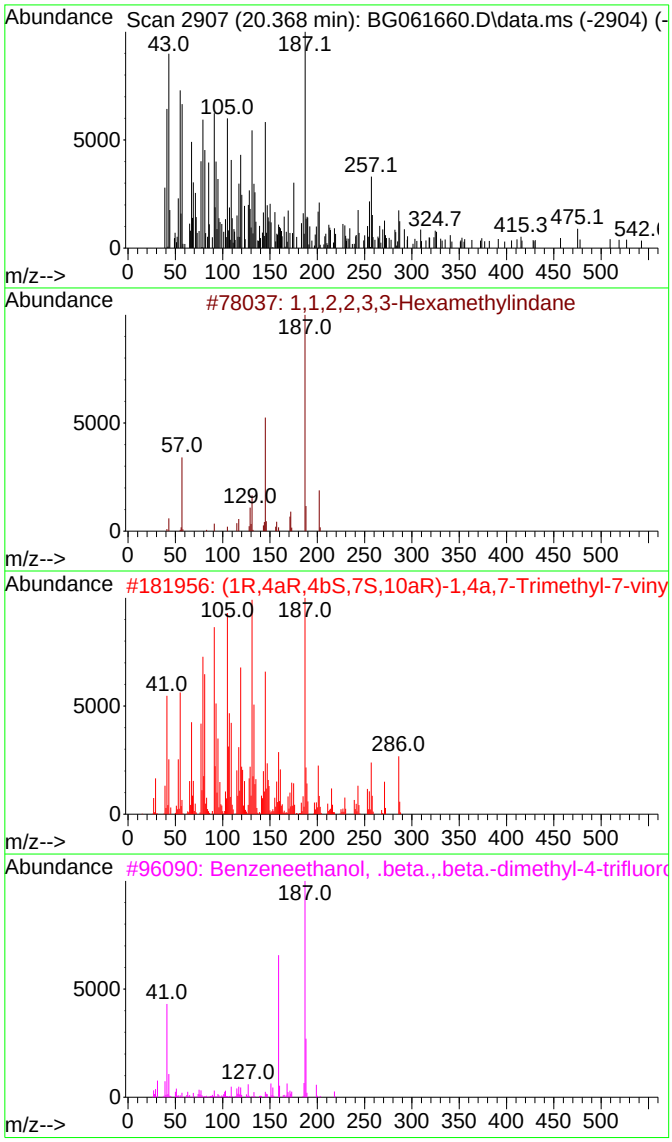
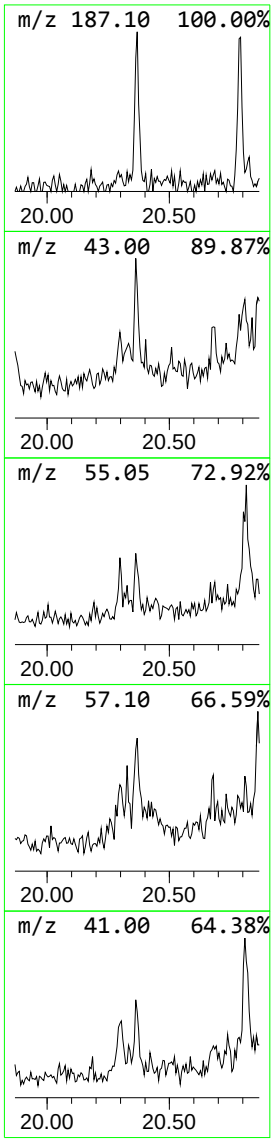
Peak Number 6 unknown-01 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.368	2.44 ng/ul	168452	Chrysene-d12	21.708

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1,2,2,3,3-Hexamethylindane	202	C15H22	091324-94-6	42
2	(1R,4aR,4bS,7S,10aR)-1,4a,7-Trim...	286	C20H30O	001686-63-1	41
3	Benzeneethanol, .beta.,.beta.-di...	218	C11H13F3O	032445-90-2	41
4	Cycloisolongifolene, 9,10-dehydro-	202	C15H22	1000156-81-6	38
5	3-Phenyl-3-trifluoromethyldiazir...	188	C8H7F3N2	040618-96-0	30





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Data File : BG061660.D
Acq On : 22 Jun 2024 22:36
Operator : MA/JU
Sample : P2776-21
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
A4C74

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG062024.MA.M
Quant Title : SVOA CALIBRATION

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

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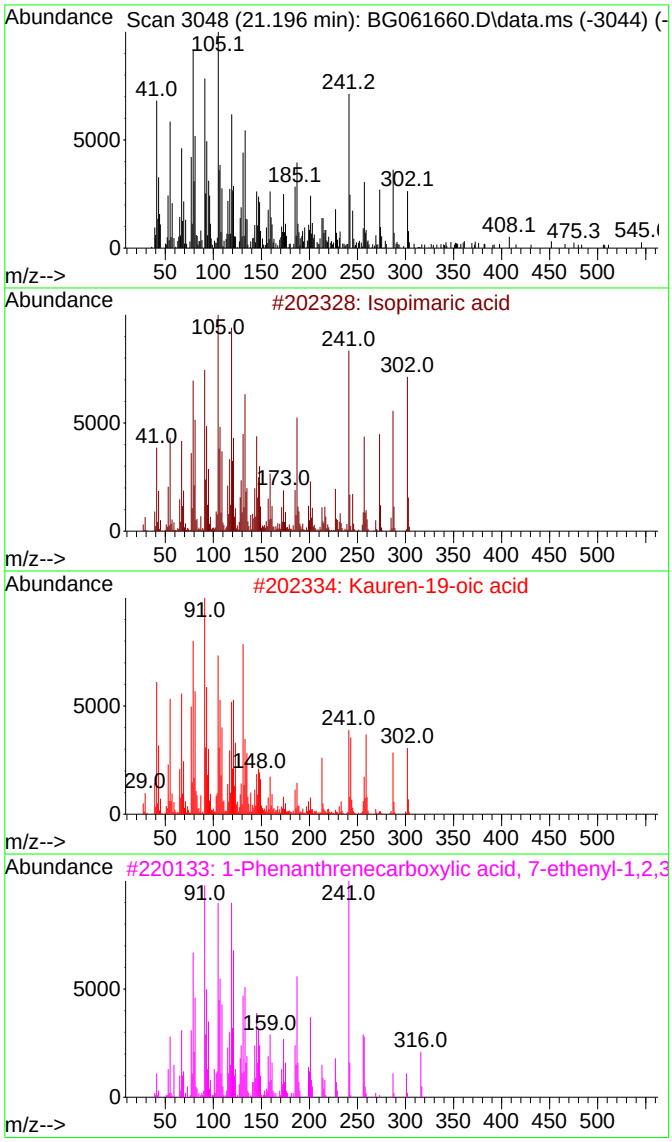
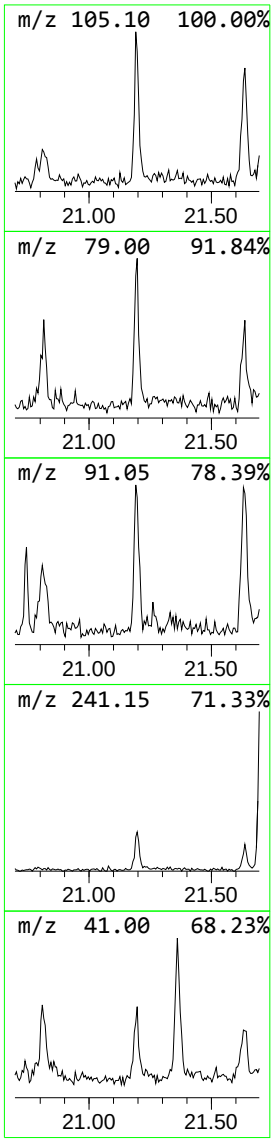
Peak Number 7 Isopimaric acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.196	4.85 ng/ul	335412	Chrysene-d12	21.708

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Isopimaric acid	302	C20H30O2	005835-26-7	97
2	Kauren-19-oic acid	302	C20H30O2	006730-83-2	60
3	1-Phenanthrenecarboxylic acid, 7...	316	C21H32O2	001686-62-0	53
4	Ethyl 5,8,11,14,17-icosapentaenoate	330	C22H34O2	084494-70-2	43
5	1-(2-Pyridinyl)-4-piperidinamine...	273	C12H14F3N3O	1488704-18-2	30





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Data File : BG061660.D
Acq On : 22 Jun 2024 22:36
Operator : MA/JU
Sample : P2776-21
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
A4C74

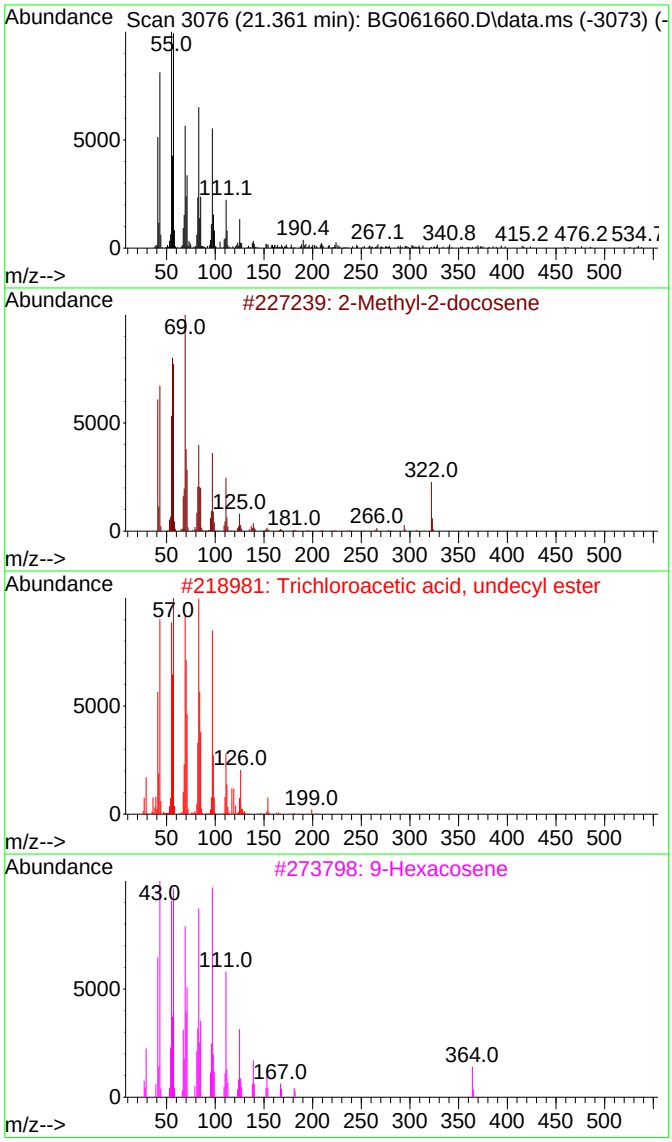
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG062024.MA.M
Quant Title : SVOA CALIBRATION

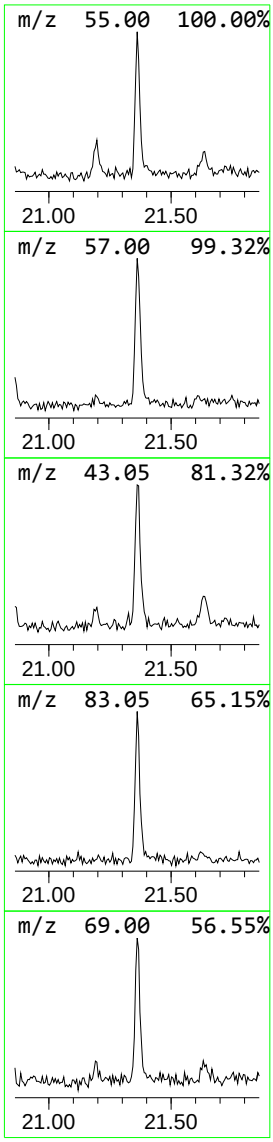
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.361	5.94 ng/ul	410403	Chrysene-d12	21.708

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Methyl-2-docosene	322	C23H46	1000131-16-9	91
2	Trichloroacetic acid, undecyl ester	316	C13H23Cl3O2	074339-49-4	87
3	9-Hexacosene	364	C26H52	071502-22-2	86
4	Cetene	224	C16H32	000629-73-2	83
5	Cetene	224	C16H32	000629-73-2	83





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062224\
Data File : BG061660.D
Acq On : 22 Jun 2024 22:36
Operator : MA/JU
Sample : P2776-21
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
A4C74

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG062024.MA.M
Quant Title : SVOA CALIBRATION

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

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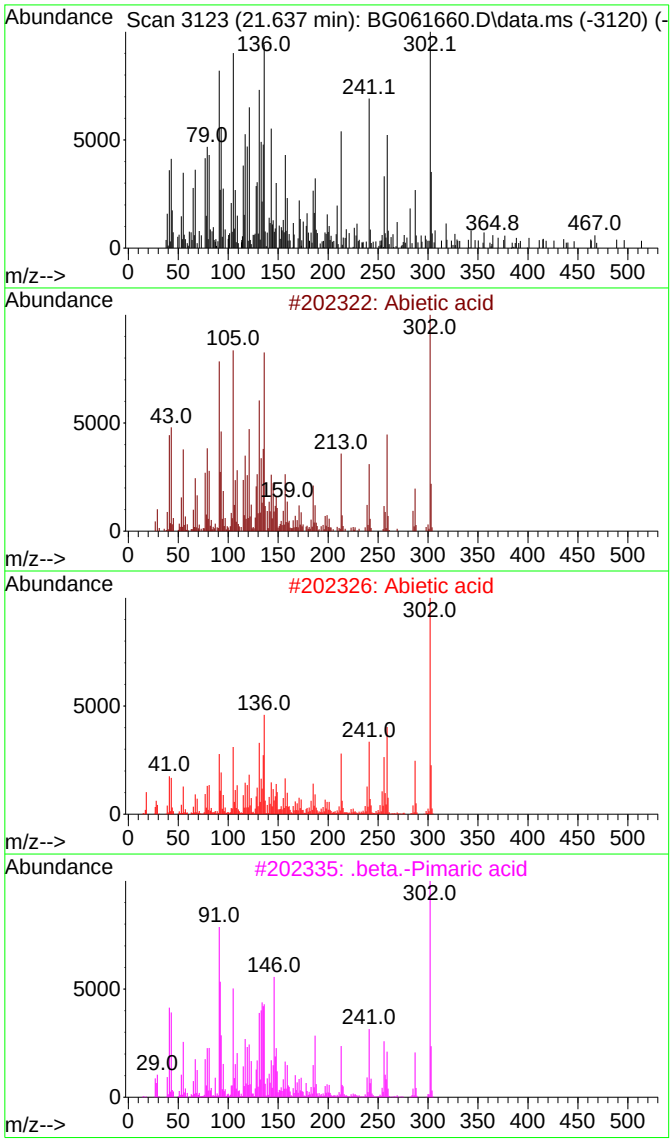
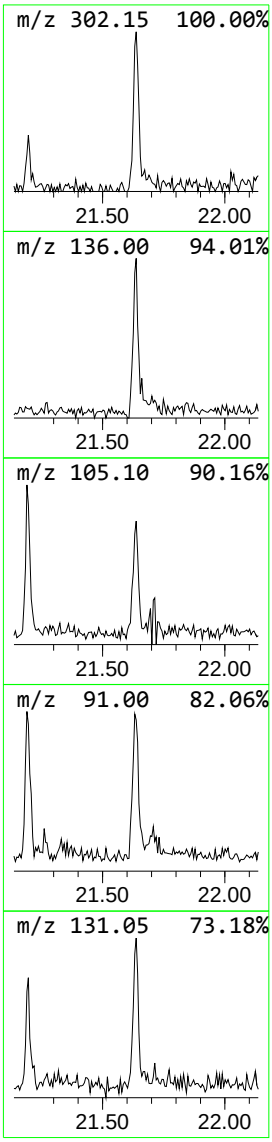
Peak Number 9 Abietic acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.637	3.96 ng/ul	273948	Chrysene-d12	21.708

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Abietic acid	302	C20H30O2	000514-10-3	93
2	Abietic acid	302	C20H30O2	000514-10-3	81
3	.beta.-Pimaric acid	302	C20H30O2	000079-54-9	70
4	Abietic acid	302	C20H30O2	000514-10-3	68
5	Abietic acid	302	C20H30O2	000514-10-3	64





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062224\
Data File : BG061660.D
Acq On : 22 Jun 2024 22:36
Operator : MA/JU
Sample : P2776-21
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
A4C74

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG062024.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
Propylene Glycol	3.740	3.3	ng/ul	85714	1	8.076	526744	20.0
Bicyclo[3.1.1]h...	6.754	18.3	ng/ul	480886	1	8.076	526744	20.0
Cyclohexene, 4-...	7.512	47.2	ng/ul	1242500	1	8.076	526744	20.0
1-Phenoxypropan...	11.625	3.2	ng/ul	140038	2	10.891	883591	20.0
n-Hexadecanoic ...	18.329	5.8	ng/ul	417711	4	17.442	1432920	20.0
unknown-01	20.368	2.4	ng/ul	168452	5	21.708	1381890	20.0
Isopimaric acid	21.196	4.8	ng/ul	335412	5	21.708	1381890	20.0
(DEL) Alkane: S...	21.361	5.9	ng/ul	410403	5	21.708	1381890	20.0
Abietic acid	21.637	4.0	ng/ul	273948	5	21.708	1381890	20.0