LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062824\
Data File : BG061779.D
Acq On : 28 Jun 2024 19:43
Operator : MA/JU
Sample : P2934-16
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
A4CQ2

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: BG061779.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.464	25	30	35	rBV4	17381	31120	1.71%	0.132%
2	3.728	70	75	82	rBV	74314	128524	7.08%	0.544%
3	3.875	96	100	113	rBV	176223	333754	18.39%	1.412%
4	3.998	115	121	124	rVV5	10963	20059	1.11%	0.085%
5	4.063	130	132	137	rVB4	5157	6492	0.36%	0.027%
6	4.216	155	158	162	rBV6	11746	13285	0.73%	0.056%
7	4.968	282	286	288	rBV4	4042	4954	0.27%	0.021%
8	5.132	308	314	319	rBV3	15425	27813	1.53%	0.118%
9	5.226	326	330	333	rBV5	5994	8936	0.49%	0.038%
10	5.367	352	354	356	rBV2	5197	4734	0.26%	0.020%
11	6.055	466	471	475	rBV5	5293	7937	0.44%	0.034%
12	6.114	475	481	489	rVB3	25361	46774	2.58%	0.198%
13	6.537	549	553	554	rVB4	3983	4605	0.25%	0.019%
14	6.607	563	565	569	rVB3	5329	5570	0.31%	0.024%
15	6.719	579	584	586	rBV3	2452	4193	0.23%	0.018%
16	6.913	614	617	621	rBV3	7937	11804	0.65%	0.050%
17	7.024	630	636	638	rBV4	6671	11331	0.62%	0.048%
18	7.206	660	667	676	rBV	363003	614265	33.84%	2.599%
19	7.383	688	697	705	rVB	235107	398327	21.94%	1.685%
20	7.535	719	723	725	rBV5	3551	6336	0.35%	0.027%
21	7.582	725	731	737	rVV	317909	537028	29.58%	2.272%
22	7.635	737	740	744	rVB2	19716	24534	1.35%	0.104%
23	7.735	754	757	760	rVB4	5677	7510	0.41%	0.032%
24	7.864	775	779	780	rBV3	4398	5021	0.28%	0.021%
25	8.058	804	812	817	rBV	338168	573501	31.59%	2.426%
26	8.105	817	820	826	rVB2	18576	28125	1.55%	0.119%
27	8.634	908	910	913	rBV3	3719	4250	0.23%	0.018%
28	8.752	921	930	939	rBV2	410330	694620	38.27%	2.939%
29	9.116	988	992	994	rBV3	5190	6274	0.35%	0.027%
30	9.139	994	996	1002	rVB4	3670	5504	0.30%	0.023%
31	9.222	1002	1010	1016	rBV	316514	571955	31.51%	2.420%
32	9.369	1034	1035	1038	rBV3	5568	5556	0.31%	0.024%
33	9.610	1075	1076	1082	rVB3	7101	9242	0.51%	0.039%
34	9.774	1098	1104	1109	rVB2	58464	94093	5.18%	0.398%
35	9.939	1125	1132	1140	rBV	266523	478976	26.39%	2.026%
36	10.209	1176	1178	1181	rVB2	4443	4644	0.26%	0.020%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062824\
Data File : BG061779.D
Acq On : 28 Jun 2024 19:43
Operator : MA/JU
Sample : P2934-16
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : A4CQ2

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	10.232	1181	1182	1186	rBV3	3374	4311	0.24%	0.018%
38	10.279	1186	1190	1194	rBV6	8558	14536	0.80%	0.061%
39	10.373	1204	1206	1212	rVB6	5499	7810	0.43%	0.033%
40	10.479	1214	1224	1234	rVV	407507	722083	39.78%	3.055%
41	10.867	1280	1290	1297	rBV	476722	869326	47.89%	3.678%
42	11.002	1306	1313	1322	rBV	119571	240510	13.25%	1.017%
43	11.390	1372	1379	1384	rBV4	29742	53470	2.95%	0.226%
44	11.601	1408	1415	1424	rVV2	86792	167904	9.25%	0.710%
45	11.748	1439	1440	1444	rVV3	4457	5173	0.28%	0.022%
46	11.801	1447	1449	1453	rVB5	3808	4270	0.24%	0.018%
47	12.018	1483	1486	1488	rVV4	4028	5643	0.31%	0.024%
48	12.071	1490	1495	1496	rVV4	3512	5660	0.31%	0.024%
49	12.112	1499	1502	1505	rVB5	3497	5256	0.29%	0.022%
50	12.259	1525	1527	1528	rBV2	5792	5232	0.29%	0.022%
51	12.406	1549	1552	1554	rBV3	4027	4628	0.25%	0.020%
52	12.441	1554	1558	1566	rVB7	7809	16937	0.93%	0.072%
53	12.553	1575	1577	1580	rVB3	4589	4305	0.24%	0.018%
54	12.847	1623	1627	1628	rBV4	4240	4942	0.27%	0.021%
55	12.882	1631	1633	1639	rVB5	6225	8845	0.49%	0.037%
56	12.941	1641	1643	1649	rVB4	5035	7293	0.40%	0.031%
57	13.141	1676	1677	1682	rBV5	3038	5062	0.28%	0.021%
58	13.211	1688	1689	1692	rBV2	4656	4290	0.24%	0.018%
59	13.499	1735	1738	1740	rBV4	6960	9298	0.51%	0.039%
60	13.599	1751	1755	1760	rBV6	8525	11799	0.65%	0.050%
61	13.852	1794	1798	1804	rVB8	6801	15798	0.87%	0.067%
62	13.916	1807	1809	1813	rBV5	3346	5129	0.28%	0.022%
63	13.993	1819	1822	1823	rBV2	5434	4417	0.24%	0.019%
64	14.087	1832	1838	1845	rVV	640621	915270	50.42%	3.872%
65	14.192	1855	1856	1858	rBV2	5686	4173	0.23%	0.018%
66	14.316	1874	1877	1882	rBV6	13741	26084	1.44%	0.110%
67	14.380	1882	1888	1898	rVB	706264	1122589	61.84%	4.749%
68	14.486	1904	1906	1909	rBV4	4588	5801	0.32%	0.025%
69	14.516	1909	1911	1914	rVV3	5494	4529	0.25%	0.019%
70	14.563	1917	1919	1920	rBV2	6160	5100	0.28%	0.022%
71	14.686	1933	1940	1947	rVB	717198	1093664	60.25%	4.627%
72	14.868	1964	1971	1980	rBV	309115	547815	30.18%	2.318%
73	15.021	1993	1997	2002	rVB7	20454	31736	1.75%	0.134%
74	15.080	2003	2007	2012	rVV8	21701	27597	1.52%	0.117%
75	15.244	2033	2035	2036	rVB2	6849	4341	0.24%	0.018%
76	15.262	2036	2038	2039	rBV2	6842	4558	0.25%	0.019%

LSC Area Percent Report

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Acq On : 28 Jun 2024 19:43
Operator : MA/JU
Sample : P2934-16
Misc :
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Instrument :
BNA_G
ClientSampleId :
A4CQ2

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

77	15.414	2062	2064	2065	rVB	9270	4551	0.25%	0.019%
78	15.514	2080	2081	2086	rVB3	11128	12531	0.69%	0.053%
79	15.673	2102	2108	2114	rBV	853714	1301404	71.69%	5.506%
80	15.779	2121	2126	2132	rVB2	165280	244732	13.48%	1.035%
81	16.161	2188	2191	2192	rBV3	8482	7897	0.44%	0.033%
82	16.313	2212	2217	2222	rVB4	23826	32830	1.81%	0.139%
83	16.384	2226	2229	2231	rBV4	14405	16768	0.92%	0.071%
84	16.478	2244	2245	2247	rBV2	6686	4348	0.24%	0.018%
85	16.813	2298	2302	2308	rVB8	45852	72781	4.01%	0.308%
86	17.430	2401	2407	2411	rBV	827842	1272028	70.08%	5.381%
87	17.471	2411	2414	2417	rVV	83941	106911	5.89%	0.452%
88	17.524	2417	2423	2430	rVV	843052	1292782	71.22%	5.469%
89	17.800	2466	2470	2473	rBV6	32240	48053	2.65%	0.203%
90	18.311	2552	2557	2562	rBV2	370153	523668	28.85%	2.215%
91	18.899	2653	2657	2662	rVB6	84917	118495	6.53%	0.501%
92	19.040	2678	2681	2685	rVB4	103913	122314	6.74%	0.517%
93	19.474	2752	2755	2760	rVB	146356	199770	11.01%	0.845%
94	19.580	2770	2773	2777	rBV3	118418	155938	8.59%	0.660%
95	19.809	2807	2812	2824	rVB2	971100	1590187	87.60%	6.727%
96	21.337	3068	3072	3078	rBV2	878703	1382817	76.18%	5.850%
97	21.689	3128	3132	3138	rVB2	620705	959923	52.88%	4.061%
98	22.524	3268	3274	3290	rVB4	704826	1815217	100.00%	7.679%
99	24.022	3524	3529	3534	rBV2	435541	838394	46.19%	3.547%
100	24.081	3535	3539	3549	rVB2	354707	764844	42.14%	3.236%

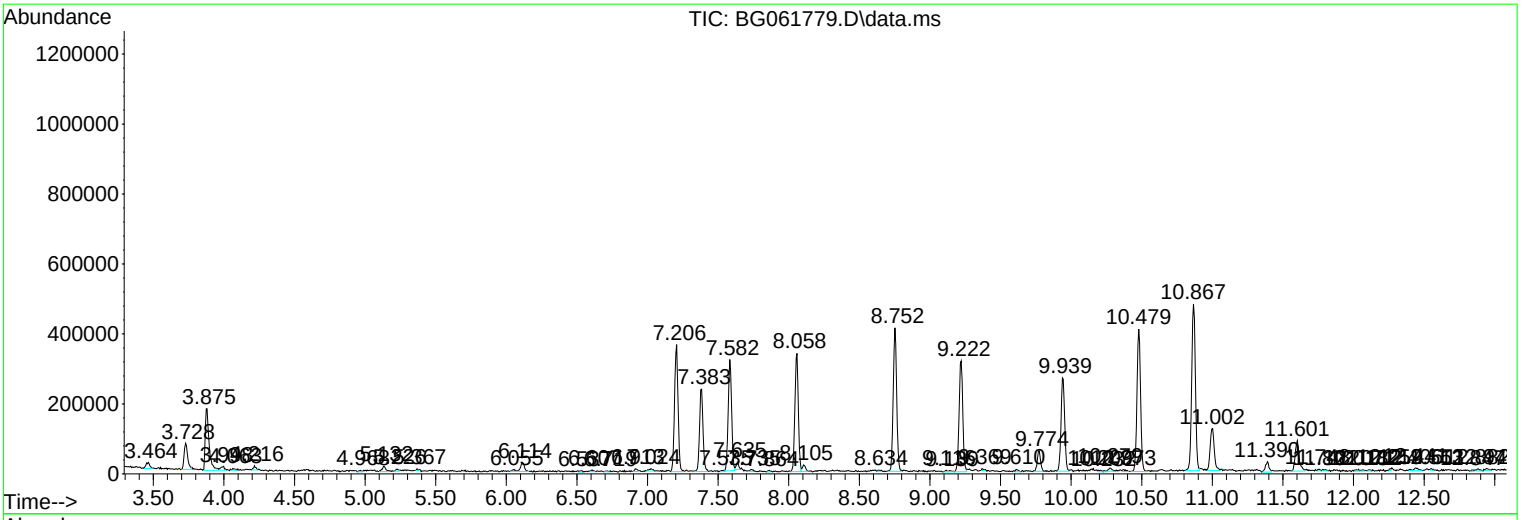
Sum of corrected areas: 23638013

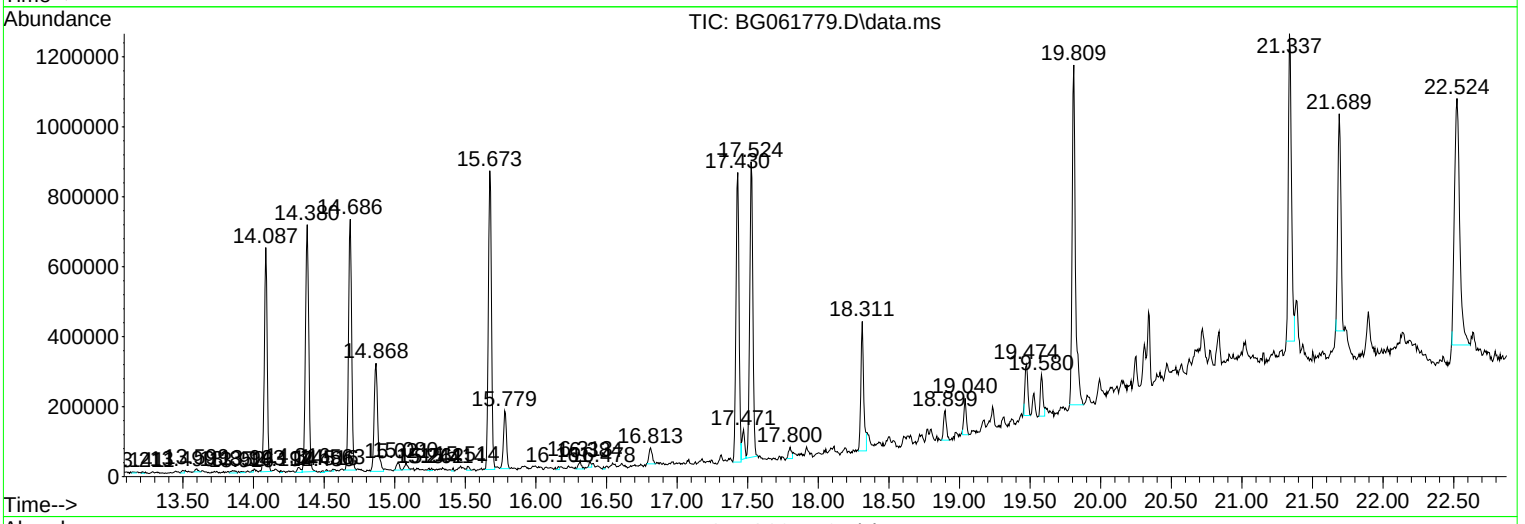
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062824\
Data File : BG061779.D
Acq On : 28 Jun 2024 19:43
Operator : MA/JU
Sample : P2934-16
Misc :
ALS Vial : 7 Sample Multiplier: 1

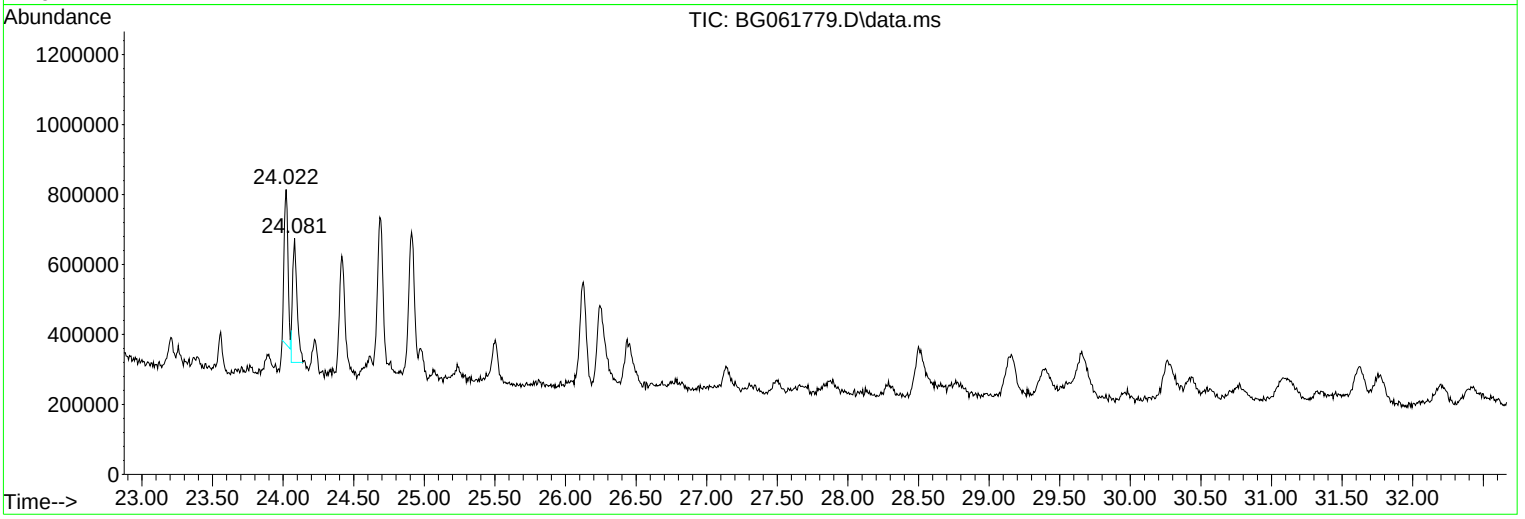
Instrument :
BNA_G
ClientSampleId :
A4CQ2

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\BG062824\
Data File : BG061779.D
Acq On : 28 Jun 2024 19:43
Operator : MA/JU
Sample : P2934-16
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
A4CQ2

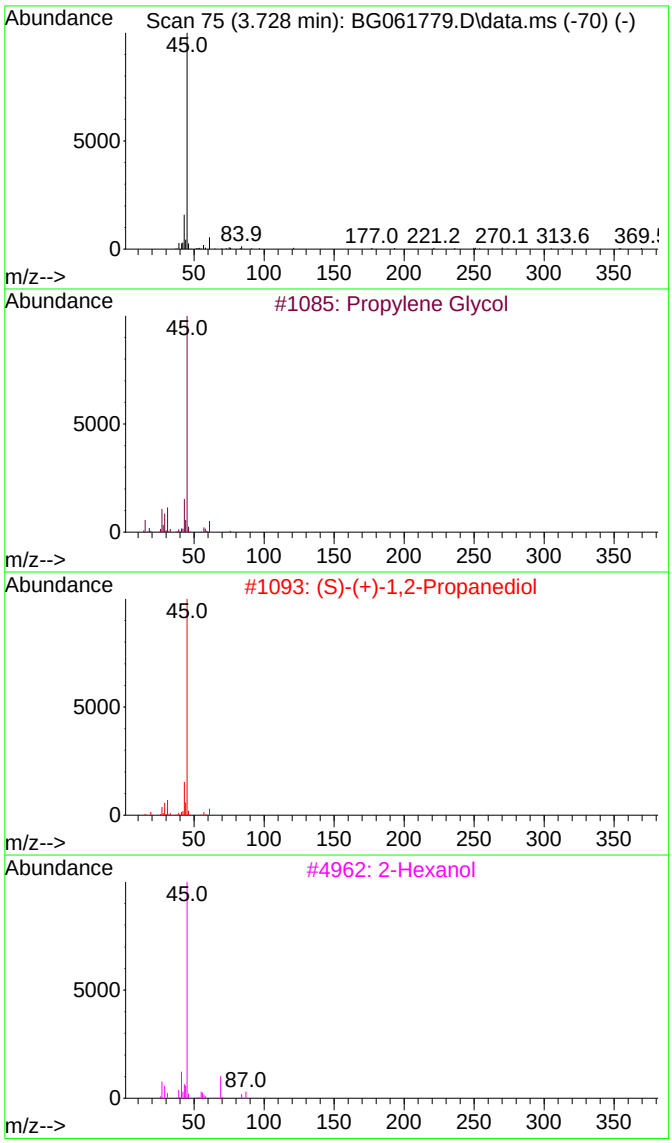
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG062024.MA.M
Quant Title : SVOA CALIBRATION

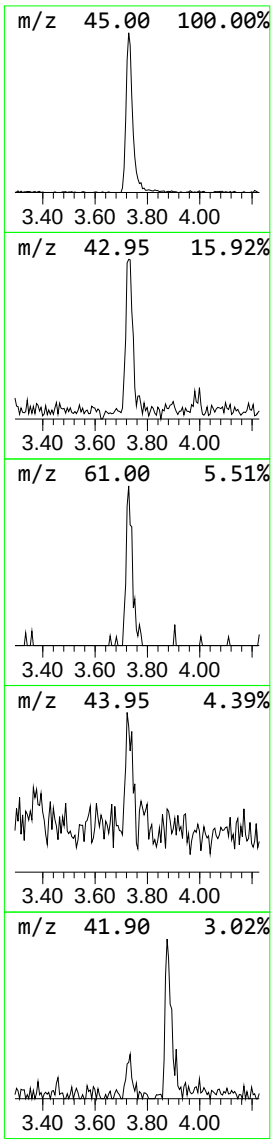
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 Propylene Glycol Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.728	4.48 ng/ul	128524	1,4-Dichlorobenzene-d4	8.058

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Propylene Glycol	76	C3H8O2	000057-55-6	50
2	(S)-(+)-1,2-Propanediol	76	C3H8O2	004254-15-3	39
3	2-Hexanol	102	C6H14O	000626-93-7	38
4	R-(-)-1,2-propanediol	76	C3H8O2	004254-14-2	25
5	Propylene Glycol	76	C3H8O2	000057-55-6	9





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062824\
Data File : BG061779.D
Acq On : 28 Jun 2024 19:43
Operator : MA/JU
Sample : P2934-16
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
A4CQ2

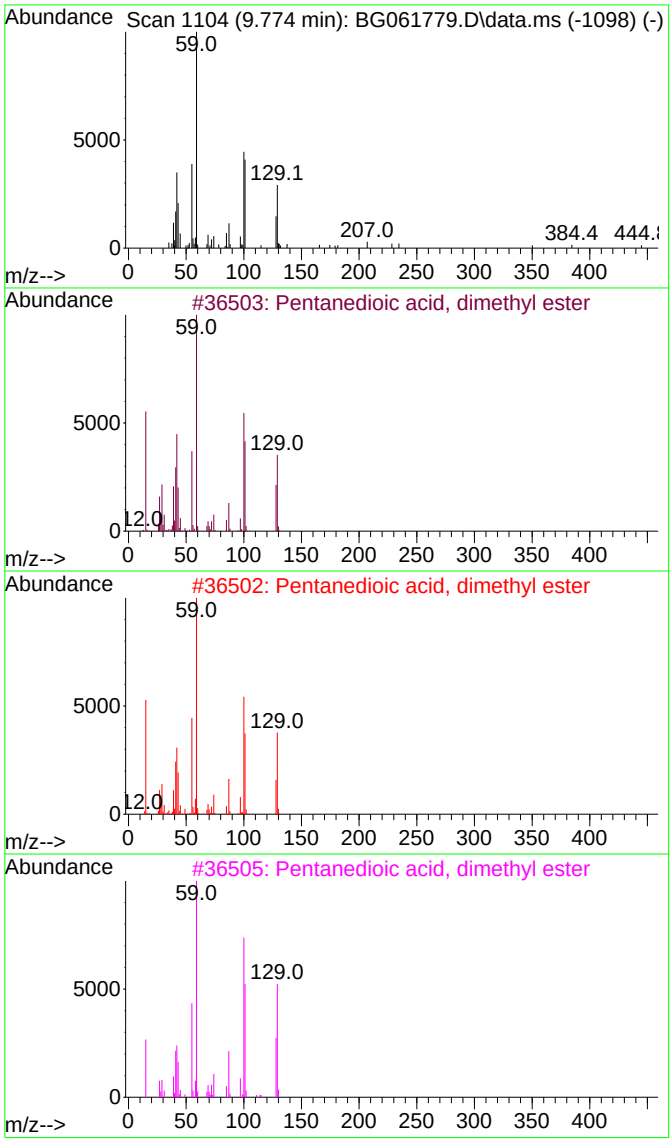
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG062024.MA.M
Quant Title : SVOA CALIBRATION

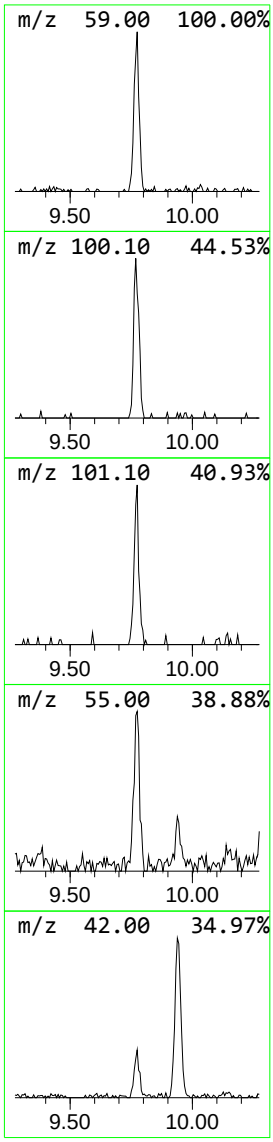
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 Pentanedioic acid, dimethyl... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.774	2.16 ng/ul	94093	Naphthalene-d8	10.867

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentanedioic acid, dimethyl ester	160	C7H12O4	001119-40-0	90
2	Pentanedioic acid, dimethyl ester	160	C7H12O4	001119-40-0	78
3	Pentanedioic acid, dimethyl ester	160	C7H12O4	001119-40-0	50
4	Pentanedioic acid, dimethyl ester	160	C7H12O4	001119-40-0	43
5	Butanedioic acid, methyl-, dimet...	160	C7H12O4	001604-11-1	40





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062824\
Data File : BG061779.D
Acq On : 28 Jun 2024 19:43
Operator : MA/JU
Sample : P2934-16
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
A4CQ2

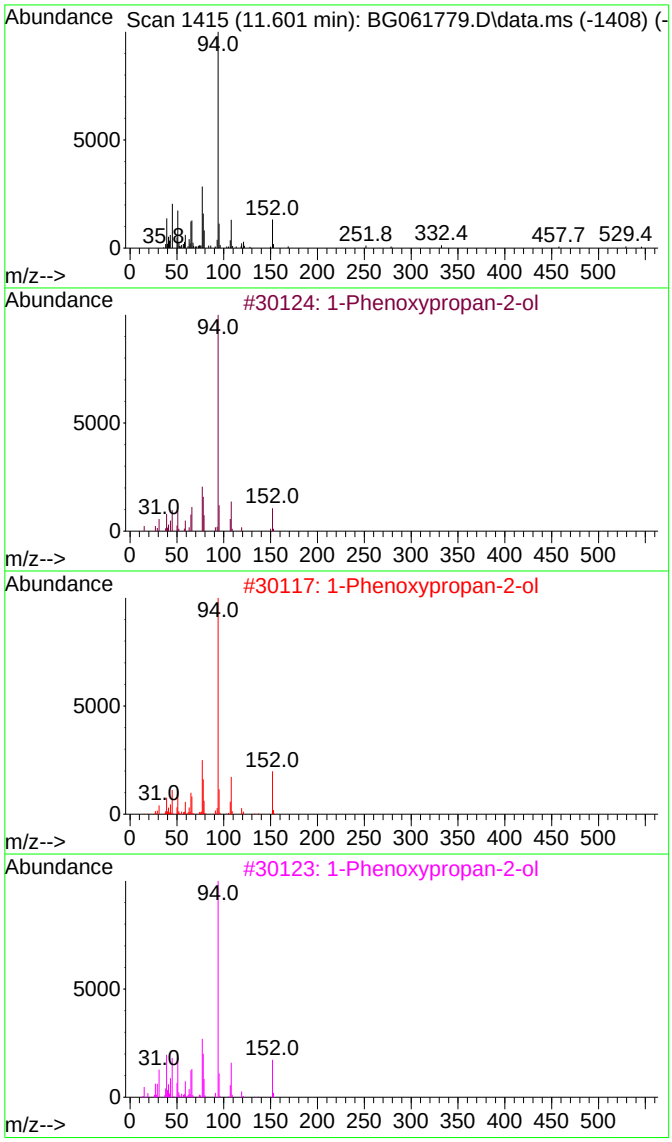
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG062024.MA.M
Quant Title : SVOA CALIBRATION

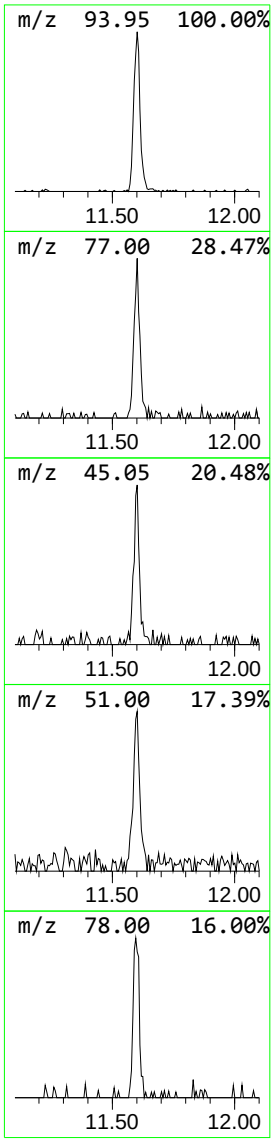
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 1-Phenoxypropan-2-ol Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.601	3.86 ng/ul	167904	Naphthalene-d8	10.867

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-Propanol, 3-phenoxy-	152	C9H12O2	006180-61-6	87
5	1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	74





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062824\
Data File : BG061779.D
Acq On : 28 Jun 2024 19:43
Operator : MA/JU
Sample : P2934-16
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
A4CQ2

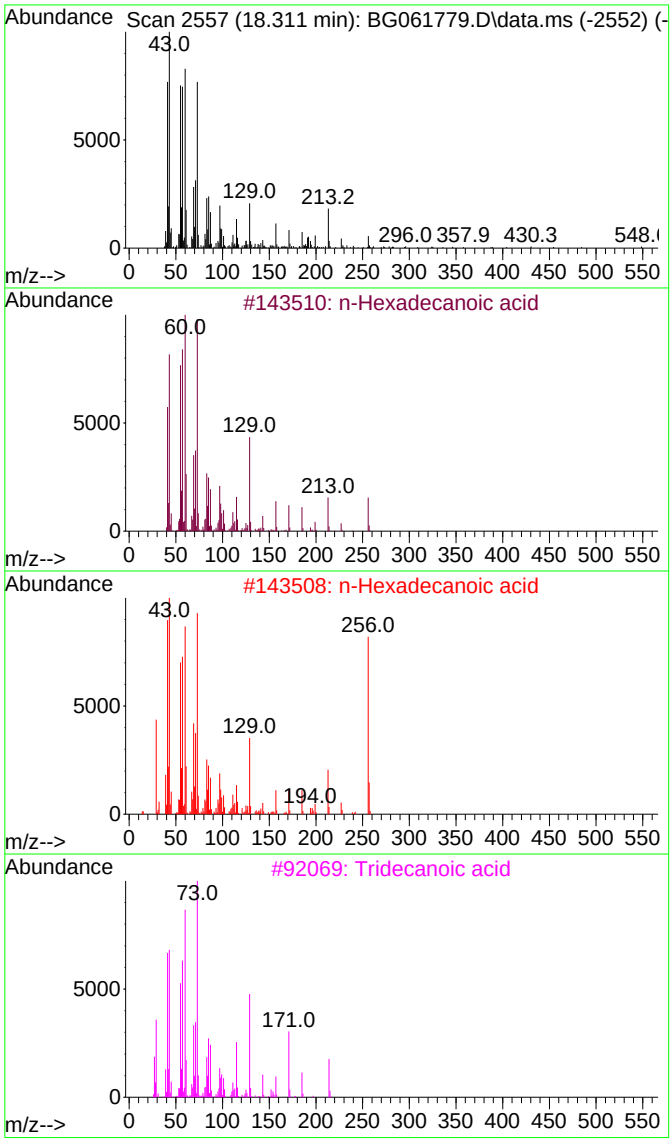
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG062024.MA.M
Quant Title : SVOA CALIBRATION

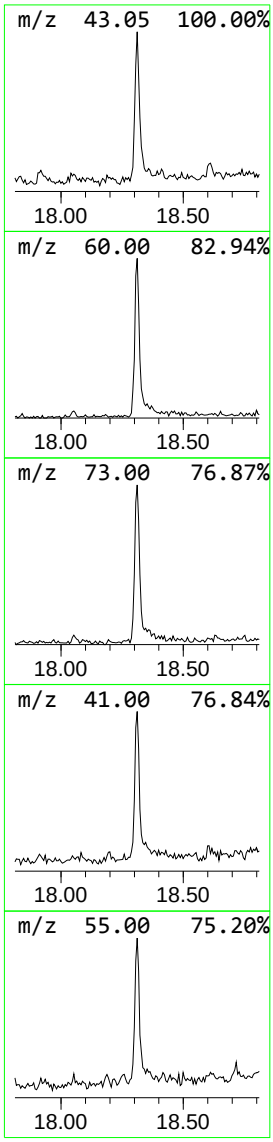
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 n-Hexadecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.311	8.23 ng/ul	523668	Phenanthrene-d10	17.430

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	97
2	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	96
3	Tridecanoic acid	214	C13H26O2	000638-53-9	89
4	Isopropyl palmitate	298	C19H38O2	000142-91-6	78
5	Tridecanoic acid	214	C13H26O2	000638-53-9	76





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062824\
Data File : BG061779.D
Acq On : 28 Jun 2024 19:43
Operator : MA/JU
Sample : P2934-16
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
A4CQ2

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

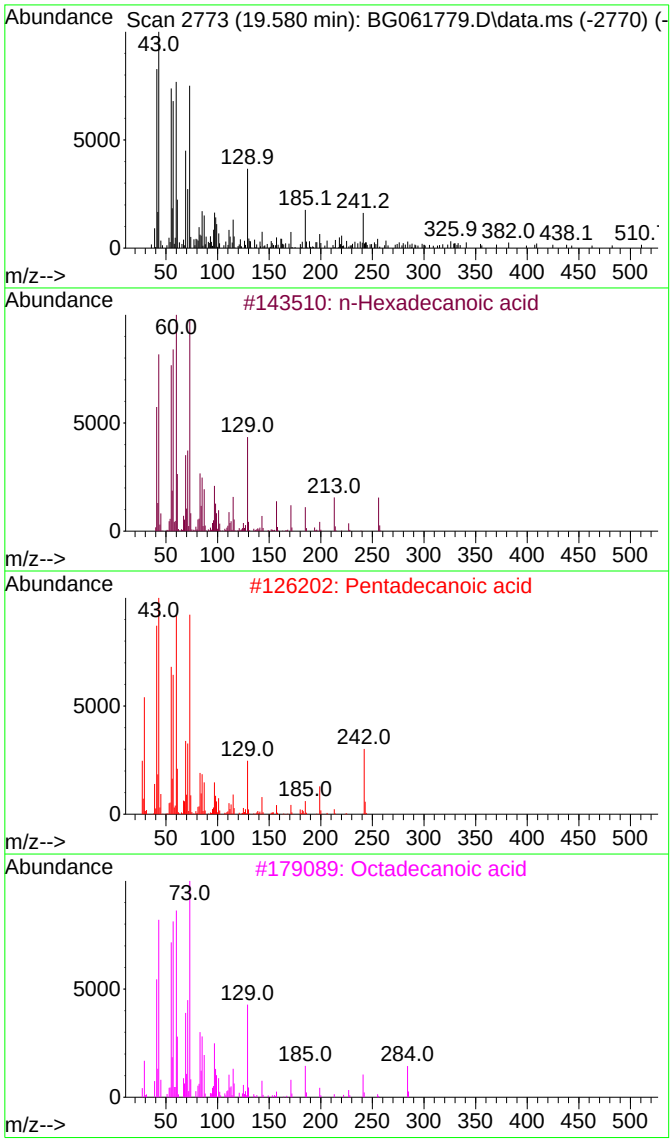
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

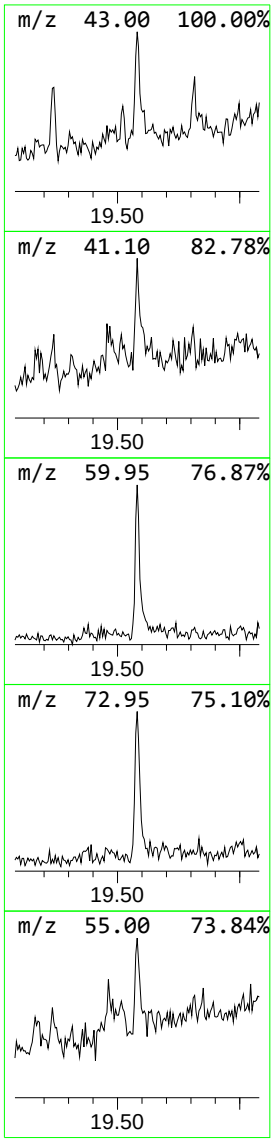
****

Peak Number 5 Pentadecanoic acid Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.580	3.25 ng/ul	155938	Chrysene-d12	21.689

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	87
2	Pentadecanoic acid	242	C15H30O2	001002-84-2	86
3	Octadecanoic acid	284	C18H36O2	000057-11-4	78
4	Pentadecanoic acid	242	C15H30O2	001002-84-2	78
5	8-Methylnonanoic acid	172	C10H20O2	005963-14-4	76





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Data File : BG061779.D
Acq On : 28 Jun 2024 19:43
Operator : MA/JU
Sample : P2934-16
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
A4CQ2

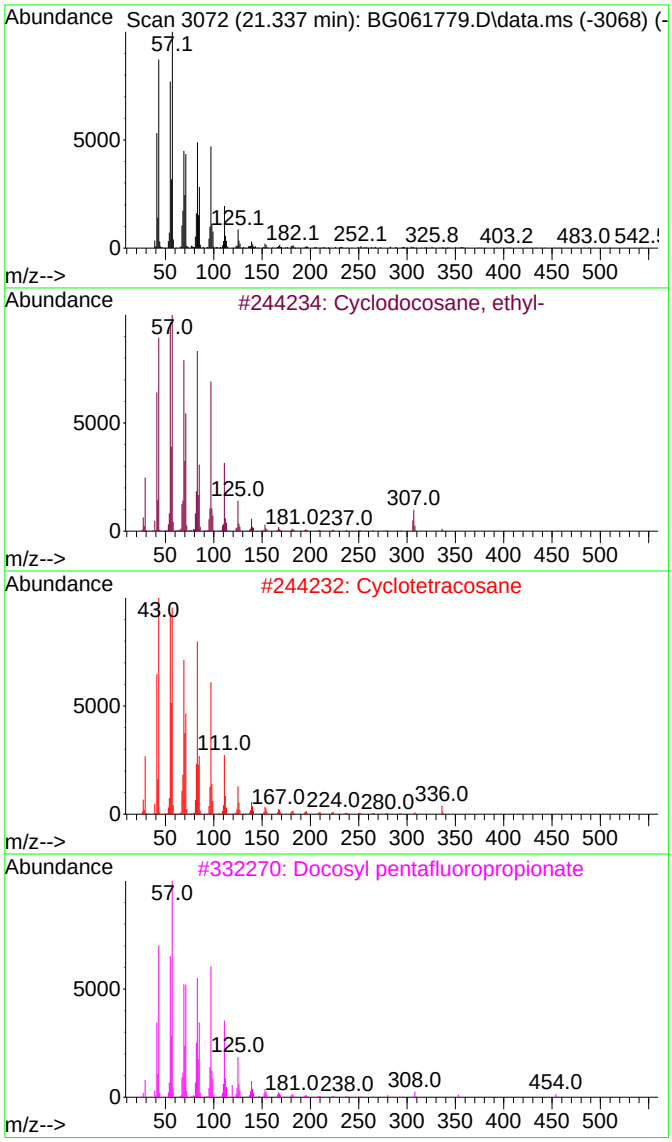
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG062024.MA.M
Quant Title : SVOA CALIBRATION

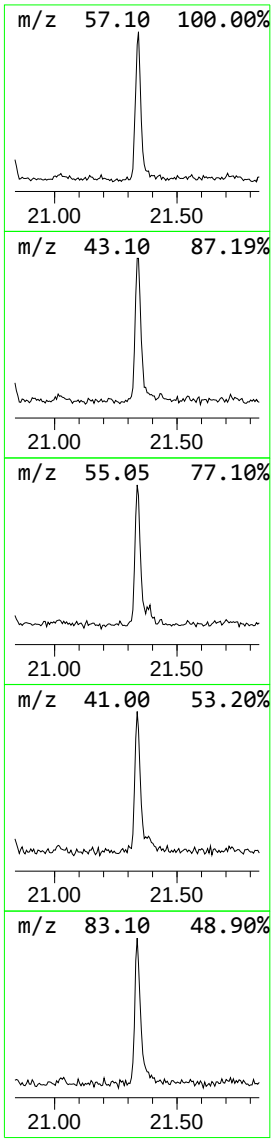
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 (DEL) Alkane: Cyclic21.337 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.337	28.81 ng/ul	1382820	Chrysene-d12	21.689

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclodocosane, ethyl-	336	C24H48	1000151-22-6	94
2	Cyclotetracosane	336	C24H48	000297-03-0	93
3	Docosyl pentafluoropropionate	472	C25H45F5O2	1000351-80-9	91
4	Tricosyl pentafluoropropionate	486	C26H47F5O2	1000351-81-0	91
5	Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	91





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Data File : BG061779.D
Acq On : 28 Jun 2024 19:43
Operator : MA/JU
Sample : P2934-16
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
A4CQ2

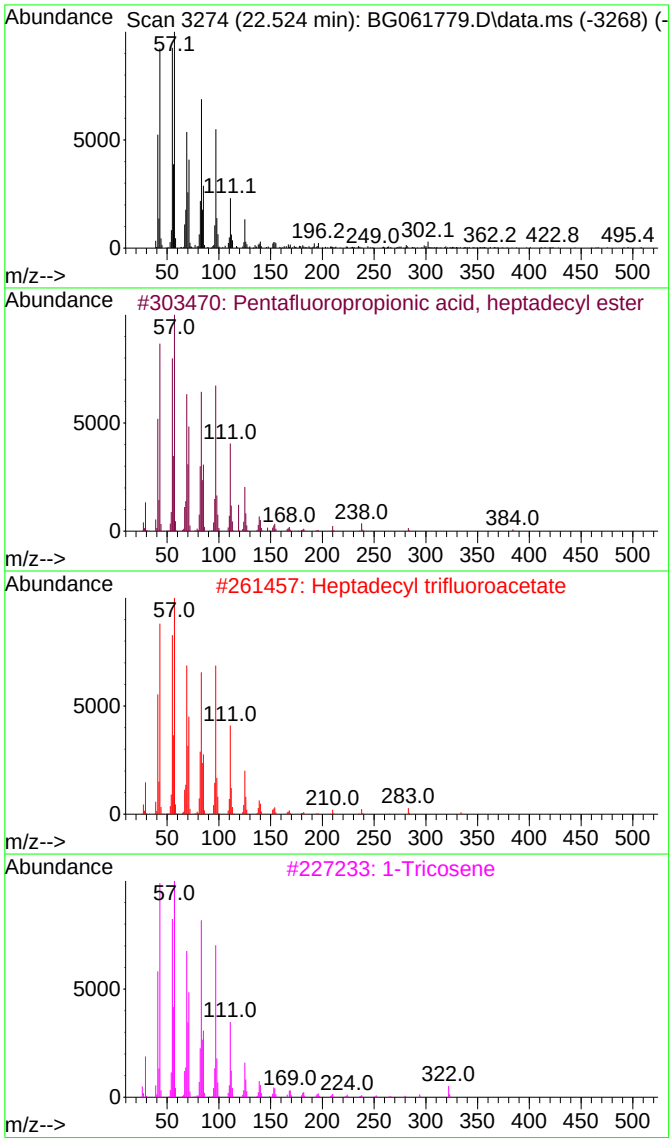
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG062024.MA.M
Quant Title : SVOA CALIBRATION

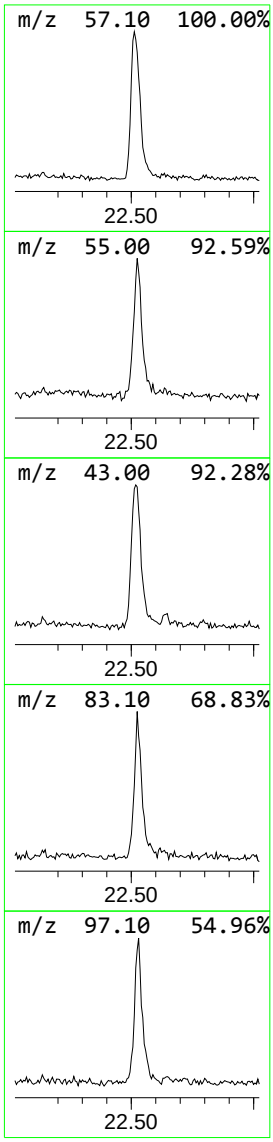
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 Pentafluoropropionic acid, ... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.524	37.82 ng/ul	1815220	Chrysene-d12	21.689

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentafluoropropionic acid, hepta...	402	C20H35F5O2	959218-78-1	91
2	Heptadecyl trifluoroacetate	352	C19H35F3O2	1010351-87-0	91
3	1-Tricosene	322	C23H46	018835-32-0	91
4	1-Docosene	308	C22H44	001599-67-3	91
5	Heptafluorobutyric acid, hexadec...	438	C20H33F7O2	006385-15-5	91





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Data File : BG061779.D
Acq On : 28 Jun 2024 19:43
Operator : MA/JU
Sample : P2934-16
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
A4CQ2

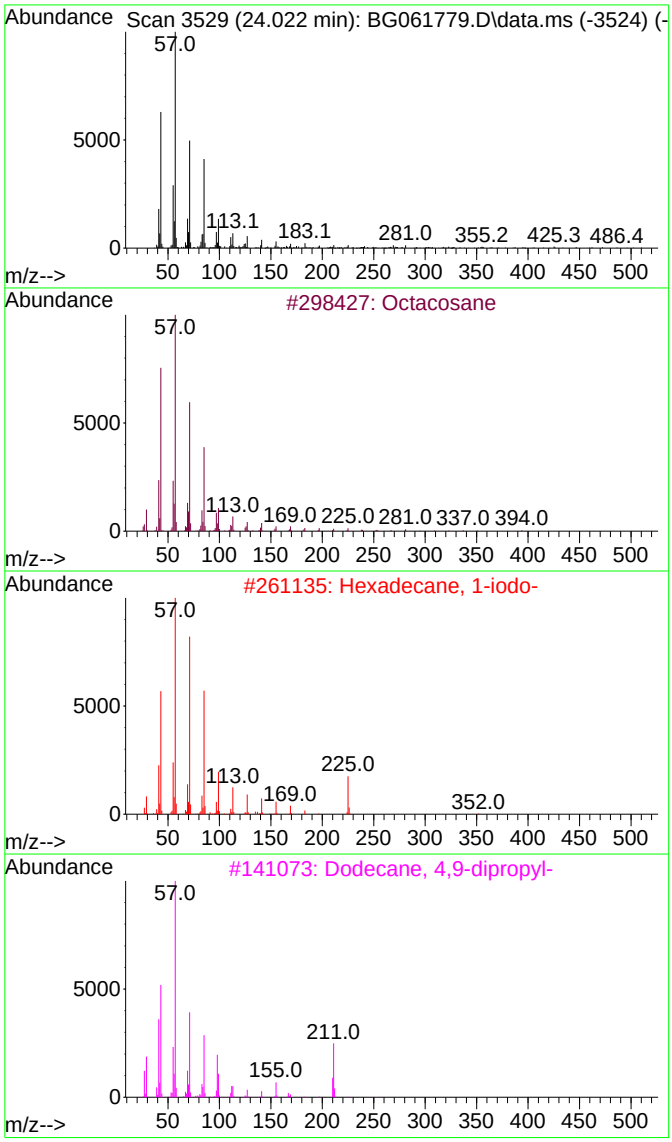
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG062024.MA.M
Quant Title : SVOA CALIBRATION

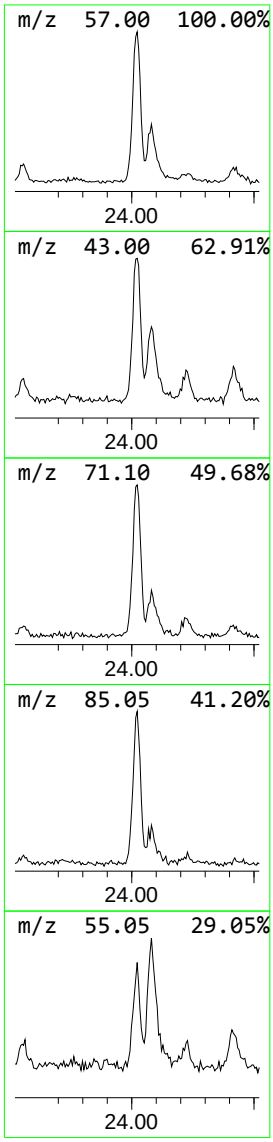
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.022	21.92 ng/ul	838394	Perylene-d12	24.909

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octacosane	394	C28H58	000630-02-4	95
2	Hexadecane, 1-iodo-	352	C16H33I	000544-77-4	94
3	Dodecane, 4,9-dipropyl-	254	C18H38	003054-63-5	87
4	Octadecane	254	C18H38	000593-45-3	83
5	Eicosane	282	C20H42	000112-95-8	83





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062824\
Data File : BG061779.D
Acq On : 28 Jun 2024 19:43
Operator : MA/JU
Sample : P2934-16
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
A4CQ2

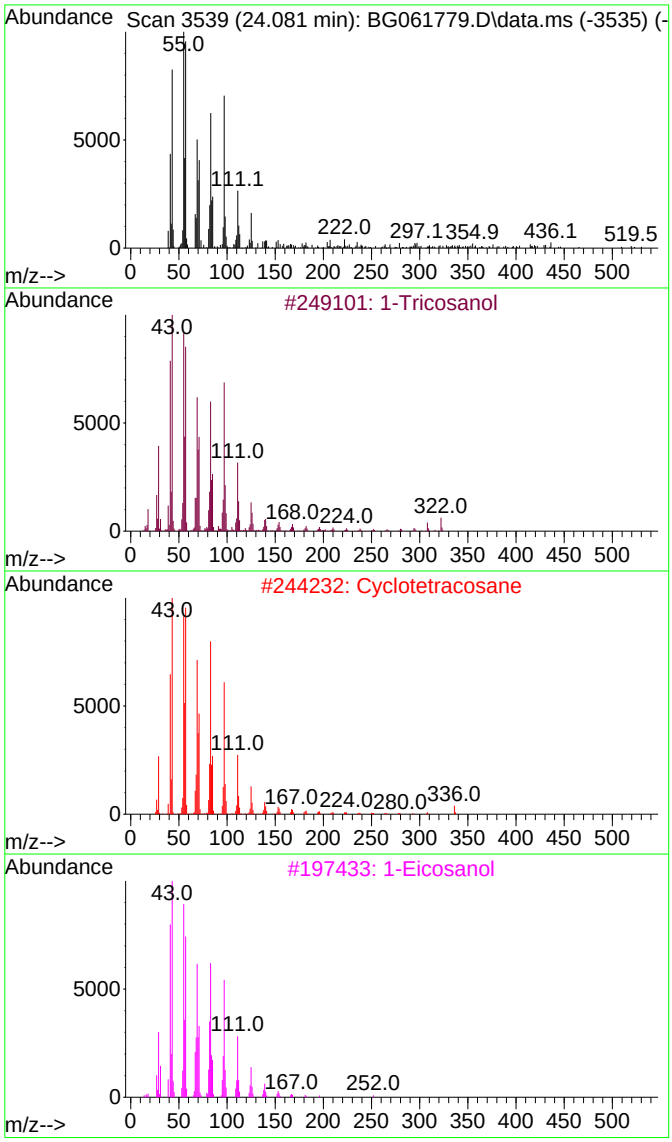
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG062024.MA.M
Quant Title : SVOA CALIBRATION

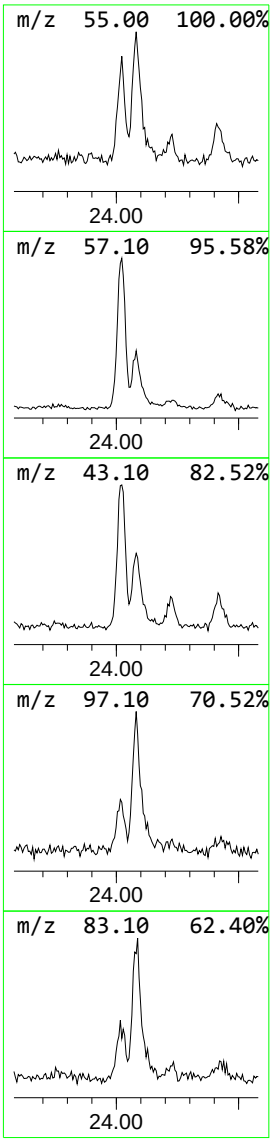
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 1-Tricosanol Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.081	20.00 ng/ul	764844	Perylene-d12	24.909

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Tricosanol	340	C23H48O	003133-01-5	91
2	Cyclotetracosane	336	C24H48	000297-03-0	91
3	1-Eicosanol	298	C20H42O	000629-96-9	91
4	Disparlure	282	C19H38O	029804-22-6	91
5	1-Docosene	308	C22H44	001599-67-3	90





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062824\
Data File : BG061779.D
Acq On : 28 Jun 2024 19:43
Operator : MA/JU
Sample : P2934-16
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
A4CQ2

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	3.728	4.5	ng/ul	128524	1	8.058	573501	20.0
Pentanedioic ac...	9.774	2.2	ng/ul	94093	2	10.867	869326	20.0
1-Phenoxypropan...	11.601	3.9	ng/ul	167904	2	10.867	869326	20.0
n-Hexadecanoic ...	18.311	8.2	ng/ul	523668	4	17.430	1272030	20.0
Pentadecanoic acid	19.580	3.3	ng/ul	155938	5	21.689	959923	20.0
(DEL) Alkane: C...	21.337	28.8	ng/ul	1382820	5	21.689	959923	20.0
Pentafluoroprop...	22.524	37.8	ng/ul	1815220	5	21.689	959923	20.0
(DEL) Alkane: S...	24.022	21.9	ng/ul	838394	6	24.909	764844	20.0
1-Tricosanol	24.081	20.0	ng/ul	764844	6	24.909	764844	20.0