LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051023\
Data File : BP015043.D
Acq On : 10 May 2023 15:28
Operator : CG/JU
Sample : 02694-18
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : JREP8

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-BP050323.MA.M
Title : SVOA CALIBRATION

Signal : TIC: BP015043.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.458	6	12	20	rVB2	83710	163727	0.83%	0.158%
2	3.740	55	60	65	rBV2	118229	160827	0.82%	0.155%
3	3.787	65	68	69	rVV	41543	44576	0.23%	0.043%
4	3.811	69	72	79	rVB	111127	146240	0.74%	0.141%
5	3.893	83	86	101	rBV	272716	412300	2.10%	0.397%
6	4.123	116	125	134	rBV	182399	319172	1.62%	0.307%
7	4.234	138	144	151	rBV2	54473	138796	0.71%	0.134%
8	4.299	153	155	164	rVB5	29597	66295	0.34%	0.064%
9	4.475	180	185	199	rVB5	33478	105596	0.54%	0.102%
10	4.611	204	208	210	rBV2	19890	24327	0.12%	0.023%
11	4.934	210	263	266	rVV2	711790	6136164	31.24%	5.906%
12	5.070	282	286	291	rBV5	17743	29623	0.15%	0.029%
13	5.193	301	307	313	rBV	234634	333058	1.70%	0.321%
14	5.434	340	348	357	rVB3	22333	59696	0.30%	0.057%
15	5.587	368	374	379	rBV	61503	96349	0.49%	0.093%
16	6.164	466	472	473	rBV3	16529	22115	0.11%	0.021%
17	6.234	473	484	491	rVV2	96334	266698	1.36%	0.257%
18	6.534	527	535	536	rBV3	14389	23149	0.12%	0.022%
19	6.564	537	540	549	rVV5	16620	49852	0.25%	0.048%
20	6.999	608	614	623	rVB	133923	208730	1.06%	0.201%
21	7.252	649	657	659	rBV3	21137	36488	0.19%	0.035%
22	7.293	659	664	673	rVB	201219	327191	1.67%	0.315%
23	7.464	687	693	708	rVB	591268	939596	4.78%	0.904%
24	7.669	720	728	734	rBV2	665614	1082131	5.51%	1.042%
25	7.728	734	738	747	rVV	99480	158110	0.80%	0.152%
26	7.922	765	771	783	rVB	214399	346024	1.76%	0.333%
27	8.146	800	809	816	rBV	607840	983036	5.00%	0.946%
28	8.205	816	819	828	rVB3	19479	36738	0.19%	0.035%
29	8.393	844	851	860	rBV	35877	69917	0.36%	0.067%
30	8.534	870	875	881	rBV	20877	35531	0.18%	0.034%
31	8.858	923	930	938	rVV	513979	851082	4.33%	0.819%
32	8.981	947	951	956	rVB2	12426	20987	0.11%	0.020%
33	9.322	1002	1009	1018	rVV	627193	1040392	5.30%	1.001%
34	9.411	1018	1024	1032	rVV5	24938	63183	0.32%	0.061%
35	9.710	1069	1075	1084	rBV2	103342	225602	1.15%	0.217%
36	9.781	1084	1087	1092	rVB6	13542	23044	0.12%	0.022%

LSC Area Percent Report

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Operator : CG/JU
Sample : 02694-18
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
JREP8

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-BP050323.MA.M
Title : SVOA CALIBRATION

37	9.881	1100	1104	1112	rVB6	10445	22254	0.11%	0.021%
38	10.052	1126	1133	1136	rBV	489821	831160	4.23%	0.800%
39	10.081	1136	1138	1144	rVV	458177	644740	3.28%	0.621%
40	10.146	1144	1149	1156	rVV	523042	843343	4.29%	0.812%
41	10.199	1156	1158	1162	rVB2	17700	23932	0.12%	0.023%
42	10.369	1178	1187	1188	rBV7	13095	25253	0.13%	0.024%
43	10.593	1216	1225	1232	rBV	798115	1327026	6.76%	1.277%
44	10.752	1247	1252	1262	rVB3	31023	65706	0.33%	0.063%
45	10.893	1270	1276	1283	rVB10	11257	26141	0.13%	0.025%
46	10.981	1283	1291	1297	rBV	825560	1376503	7.01%	1.325%
47	11.116	1306	1314	1328	rBV	621454	1132179	5.76%	1.090%
48	11.505	1375	1380	1386	rBV	57646	99061	0.50%	0.095%
49	11.569	1386	1391	1395	rVV	60752	103920	0.53%	0.100%
50	11.640	1397	1403	1413	rVB2	38242	88322	0.45%	0.085%
51	11.752	1413	1422	1432	rBV	105675	247451	1.26%	0.238%
52	11.987	1458	1462	1468	rVB5	15903	25411	0.13%	0.024%
53	12.081	1471	1478	1484	rBV7	13614	34513	0.18%	0.033%
54	12.216	1489	1501	1512	rBV	490626	915483	4.66%	0.881%
55	12.681	1555	1580	1608	rBV	4647222	19643530	100.00%	18.906%
56	13.246	1672	1676	1686	rVB2	96812	174868	0.89%	0.168%
57	13.381	1695	1699	1703	rBV7	13833	23123	0.12%	0.022%
58	13.575	1723	1732	1742	rBV	402700	694314	3.53%	0.668%
59	13.675	1745	1749	1754	rVV2	37950	74626	0.38%	0.072%
60	13.740	1754	1760	1766	rVV	488080	723340	3.68%	0.696%
61	13.804	1768	1771	1781	rVB2	65226	136994	0.70%	0.132%
62	14.022	1805	1808	1816	rVB4	45972	72496	0.37%	0.070%
63	14.099	1817	1821	1825	rVB5	19436	30633	0.16%	0.029%
64	14.187	1830	1836	1842	rBV	1514483	2128681	10.84%	2.049%
65	14.322	1854	1859	1870	rBV	339931	523731	2.67%	0.504%
66	14.481	1880	1886	1895	rVB	1716803	2622903	13.35%	2.524%
67	14.587	1900	1904	1909	rVB2	53991	81573	0.42%	0.079%
68	14.675	1916	1919	1926	rVB3	50955	76034	0.39%	0.073%
69	14.787	1932	1938	1945	rVB	1244428	1851023	9.42%	1.782%
70	14.975	1965	1970	1979	rBV	146267	261760	1.33%	0.252%
71	15.193	2003	2007	2012	rVB3	38401	63627	0.32%	0.061%
72	15.287	2018	2023	2029	rBV	244628	374889	1.91%	0.361%
73	15.593	2072	2075	2080	rBV7	29798	37081	0.19%	0.036%
74	15.781	2100	2107	2115	rBV	2236961	3285201	16.72%	3.162%
75	15.893	2121	2126	2137	rBV	441478	608891	3.10%	0.586%
76	16.098	2158	2161	2163	rBV2	24034	32882	0.17%	0.032%

LSC Area Percent Report

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Acq On : 10 May 2023 15:28
Operator : CG/JU
Sample : 02694-18
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
JREP8

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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Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050323.MA.M
Title : SVOA CALIBRATION

77	16.140	2163	2168	2173	rVB	850694	1101052	5.61%	1.060%
78	16.457	2220	2222	2227	rVB4	21033	25668	0.13%	0.025%
79	16.834	2281	2286	2297	rBV2	240801	590975	3.01%	0.569%
80	16.922	2297	2301	2308	rVB3	67618	115574	0.59%	0.111%
81	17.045	2311	2322	2340	rBV	7097867	15150728	77.13%	14.582%
82	17.287	2360	2363	2368	rVB2	82759	109814	0.56%	0.106%
83	17.540	2401	2406	2417	rBV	1628740	2634060	13.41%	2.535%
84	17.639	2418	2423	2430	rVV2	2753600	3903713	19.87%	3.757%
85	17.739	2436	2440	2451	rVB	202688	345736	1.76%	0.333%
86	18.287	2528	2533	2541	rBV	200044	373146	1.90%	0.359%
87	18.404	2547	2553	2562	rBV	1464906	2231492	11.36%	2.148%
88	18.528	2570	2574	2590	rVB	211836	347176	1.77%	0.334%
89	18.804	2618	2621	2628	rVB2	88042	120765	0.61%	0.116%
90	19.098	2668	2671	2679	rVB5	100358	141520	0.72%	0.136%
91	19.222	2689	2692	2698	rBV2	82558	121333	0.62%	0.117%
92	19.510	2738	2741	2743	rBV2	109398	141832	0.72%	0.137%
93	19.622	2756	2760	2770	rBV	799307	1302125	6.63%	1.253%
94	19.857	2794	2800	2816	rVB2	3648454	6424453	32.71%	6.183%
95	21.292	3040	3044	3052	rBV	631600	877517	4.47%	0.845%
96	21.622	3095	3100	3108	rBV	2097172	2759122	14.05%	2.656%
97	22.927	3318	3322	3331	rVB	406578	627322	3.19%	0.604%
98	24.033	3503	3510	3523	rVB2	2365188	5046268	25.69%	4.857%
99	24.204	3532	3539	3550	rVB	1368864	3006170	15.30%	2.893%

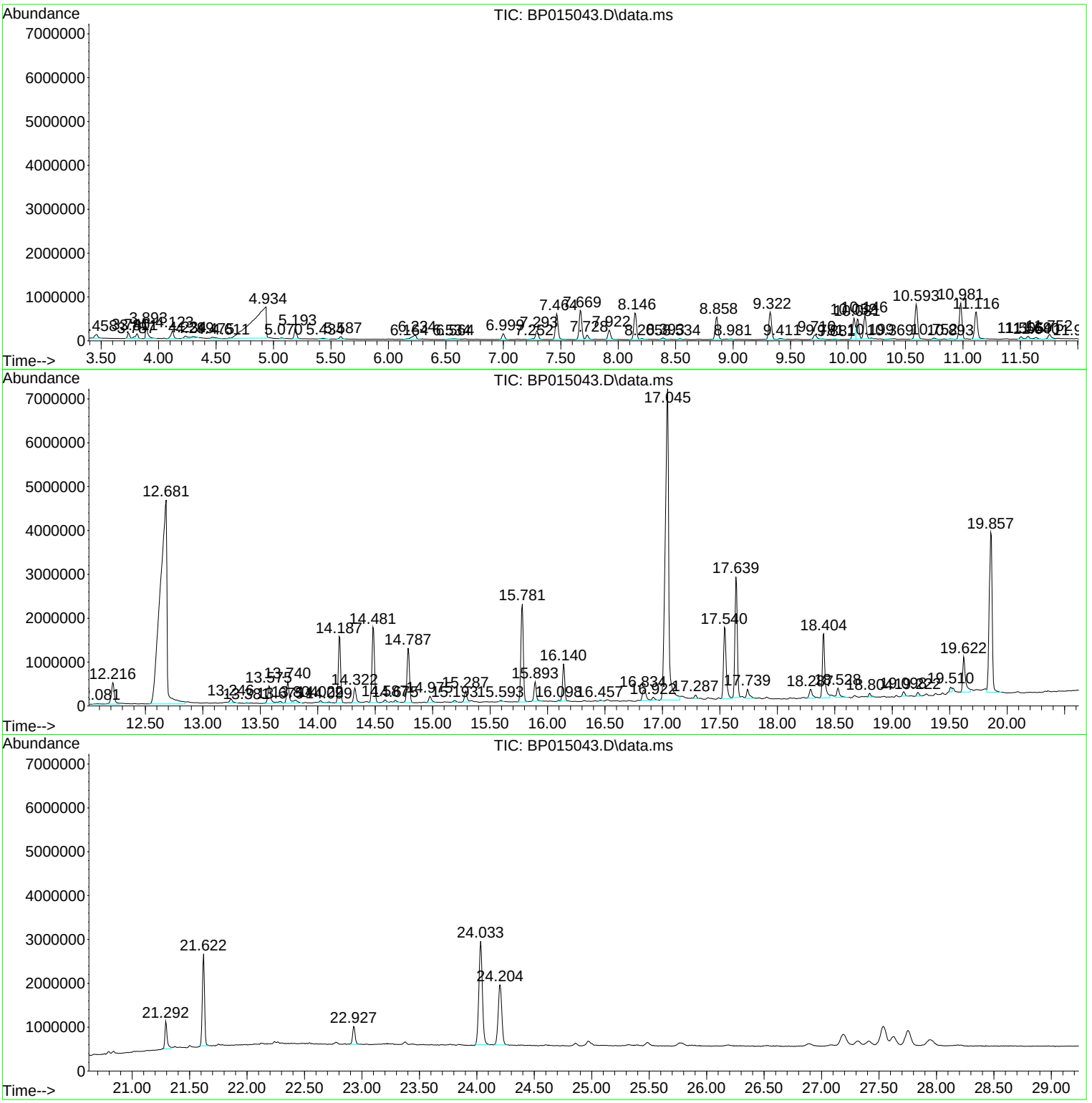
Sum of corrected areas: 103898501

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051023\
Data File : BP015043.D
Acq On : 10 May 2023 15:28
Operator : CG/JU
Sample : 02694-18
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
JREP8

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050323.MA.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051023\
Data File : BP015043.D
Acq On : 10 May 2023 15:28
Operator : CG/JU
Sample : 02694-18
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
JREP8

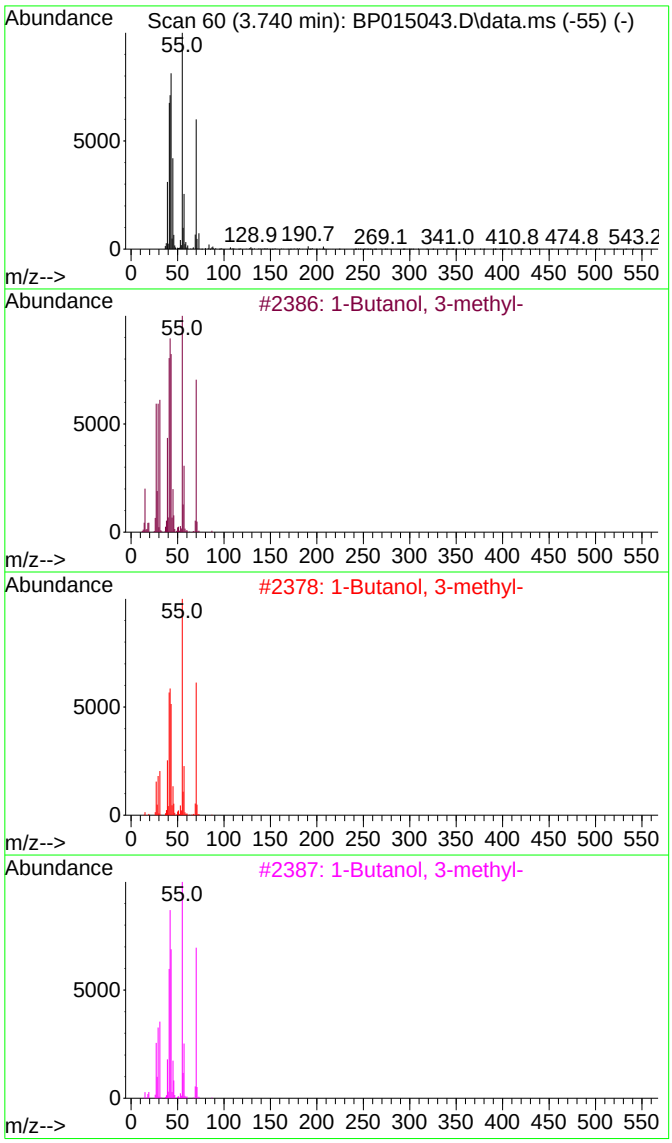
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050323.MA.M
Quant Title : SVOA CALIBRATION

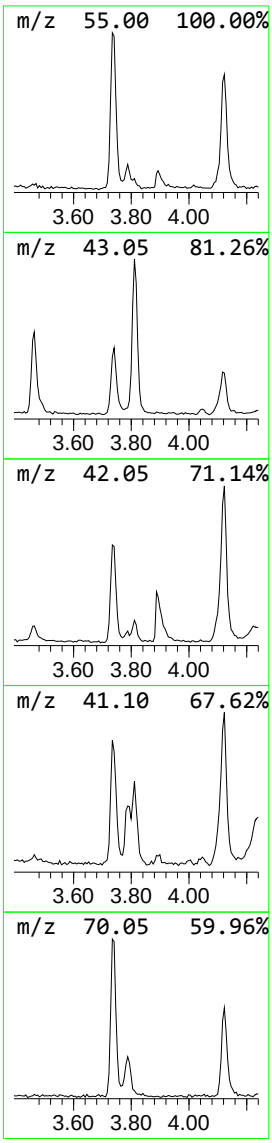
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 1-Butanol, 3-methyl- Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.740	3.27 ng/ul	160827	1,4-Dichlorobenzene-d4	8.146

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Butanol, 3-methyl-	88	C5H12O	000123-51-3	64
2	1-Butanol, 3-methyl-	88	C5H12O	000123-51-3	64
3	1-Butanol, 3-methyl-	88	C5H12O	000123-51-3	64
4	3,3-Dimethyl-1,2-epoxybutane	100	C6H12O	002245-30-9	47
5	Pentane, 1-chloro-	106	C5H11Cl	000543-59-9	47





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051023\
Data File : BP015043.D
Acq On : 10 May 2023 15:28
Operator : CG/JU
Sample : 02694-18
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
JREP8

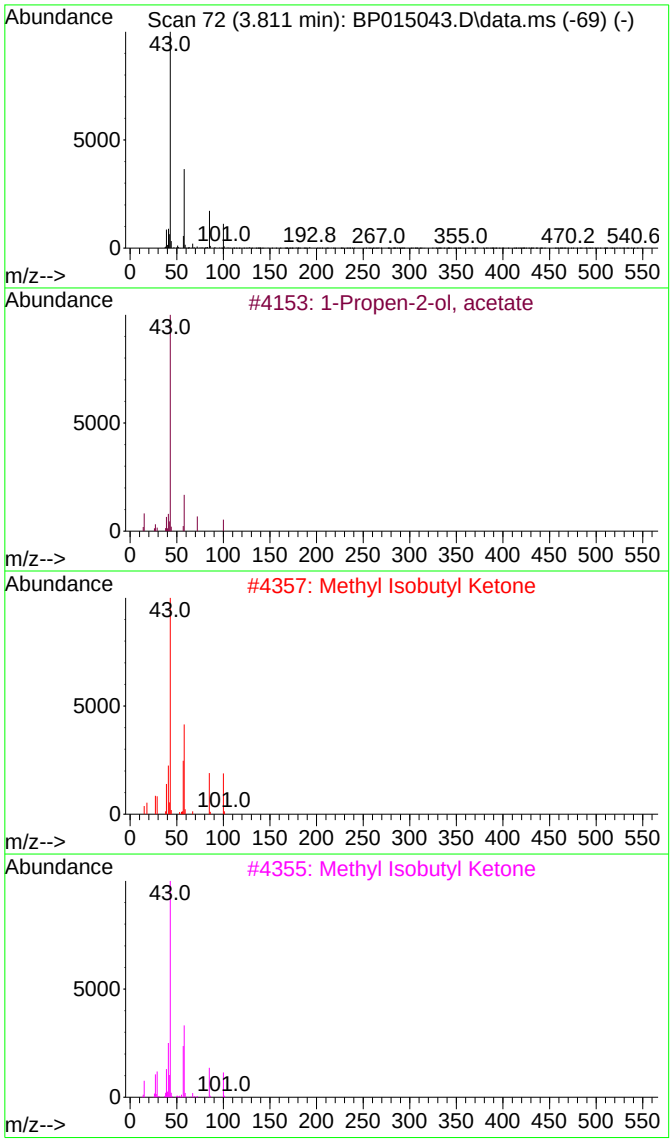
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050323.MA.M
Quant Title : SVOA CALIBRATION

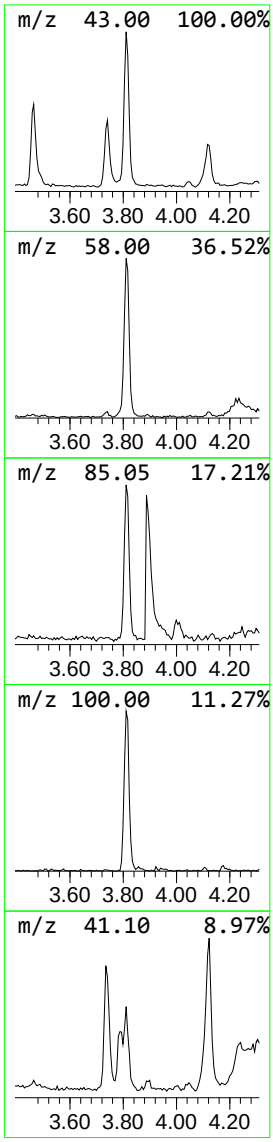
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 1-Propen-2-ol, acetate Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.811	2.98 ng/ul	146240	1,4-Dichlorobenzene-d4	8.146

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	64
2	Methyl Isobutyl Ketone	100	C6H12O	000108-10-1	59
3	Methyl Isobutyl Ketone	100	C6H12O	000108-10-1	59
4	Methyl Isobutyl Ketone	100	C6H12O	000108-10-1	42
5	2-Heptanone, 4-methyl-	128	C8H16O	006137-06-0	38





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051023\
Data File : BP015043.D
Acq On : 10 May 2023 15:28
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Sample : 02694-18
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
JREP8

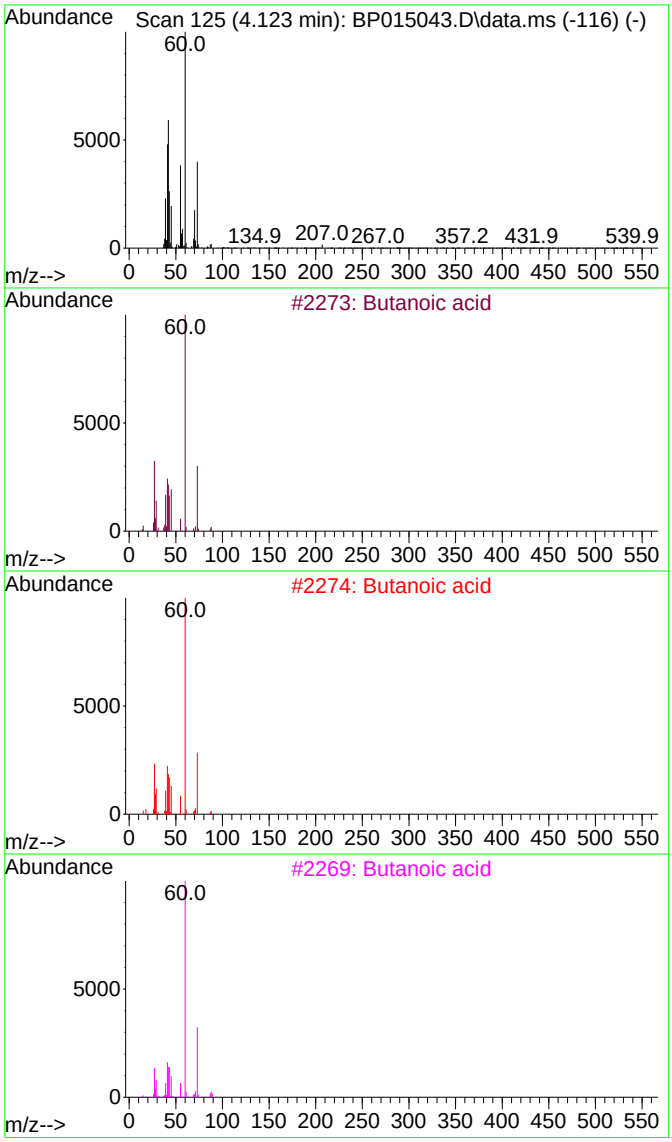
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050323.MA.M
Quant Title : SVOA CALIBRATION

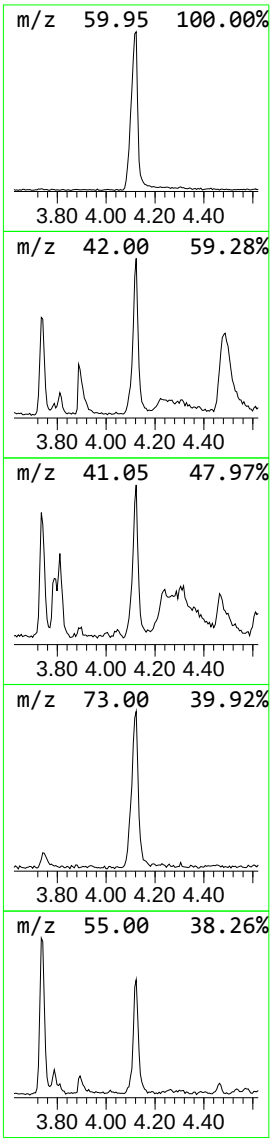
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Butanoic acid Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.123	6.49 ng/ul	319172	1,4-Dichlorobenzene-d4	8.146

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butanoic acid	88	C4H8O2	000107-92-6	72
2	Butanoic acid	88	C4H8O2	000107-92-6	64
3	Butanoic acid	88	C4H8O2	000107-92-6	59
4	Butanoic acid	88	C4H8O2	000107-92-6	53
5	Pentanoic acid	102	C5H10O2	000109-52-4	40





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051023\
Data File : BP015043.D
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Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
JREP8

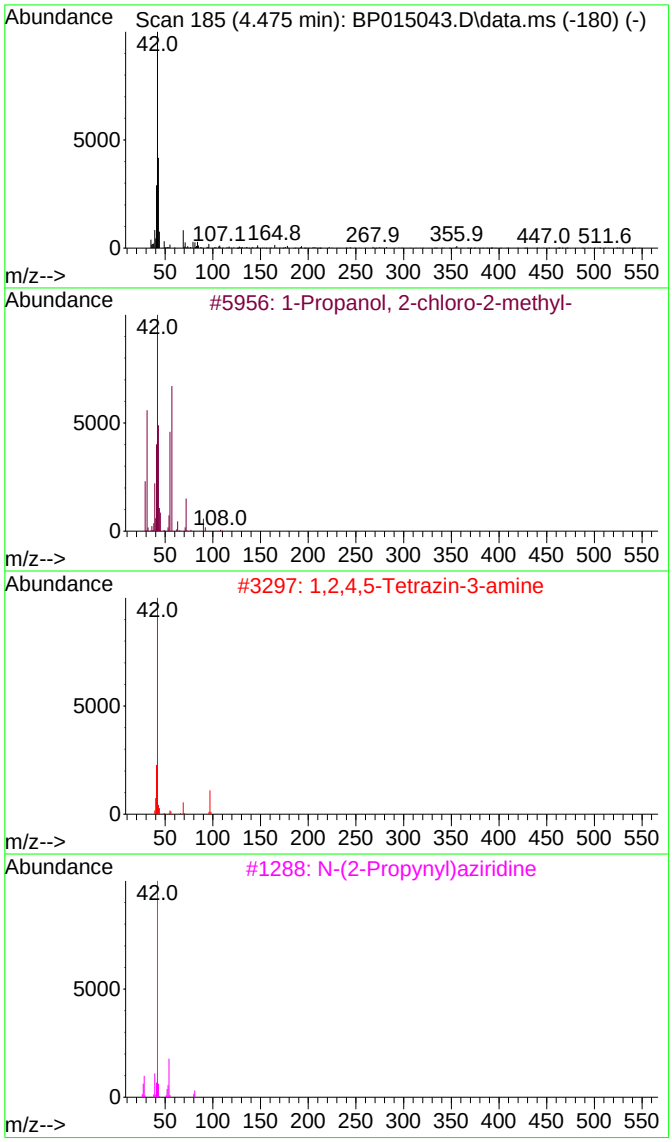
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050323.MA.M
Quant Title : SVOA CALIBRATION

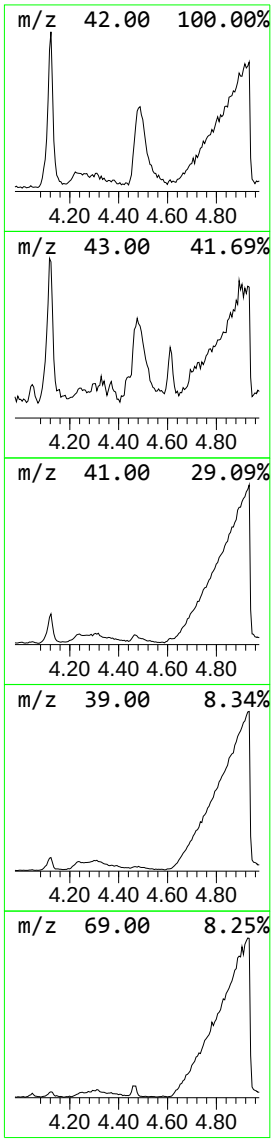
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 unknown-01 Concentration Rank 30

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.475	2.15 ng/ul	105596	1,4-Dichlorobenzene-d4	8.146

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Propanol, 2-chloro-2-methyl-	108	C4H9ClO	000558-38-3	9
2	1,2,4,5-Tetrazin-3-amine	97	C2H3N5	079329-74-1	9
3	N-(2-Propynyl)aziridine	81	C5H7N	041104-29-4	7
4	1,4-Dioxaspiro[2.4]heptan-5-one,...	128	C6H8O3	1010143-02-1	7
5	Ethylenimine	43	C2H5N	000151-56-4	5





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051023\
Data File : BP015043.D
Acq On : 10 May 2023 15:28
Operator : CG/JU
Sample : 02694-18
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
JREP8

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050323.MA.M
Quant Title : SVOA CALIBRATION

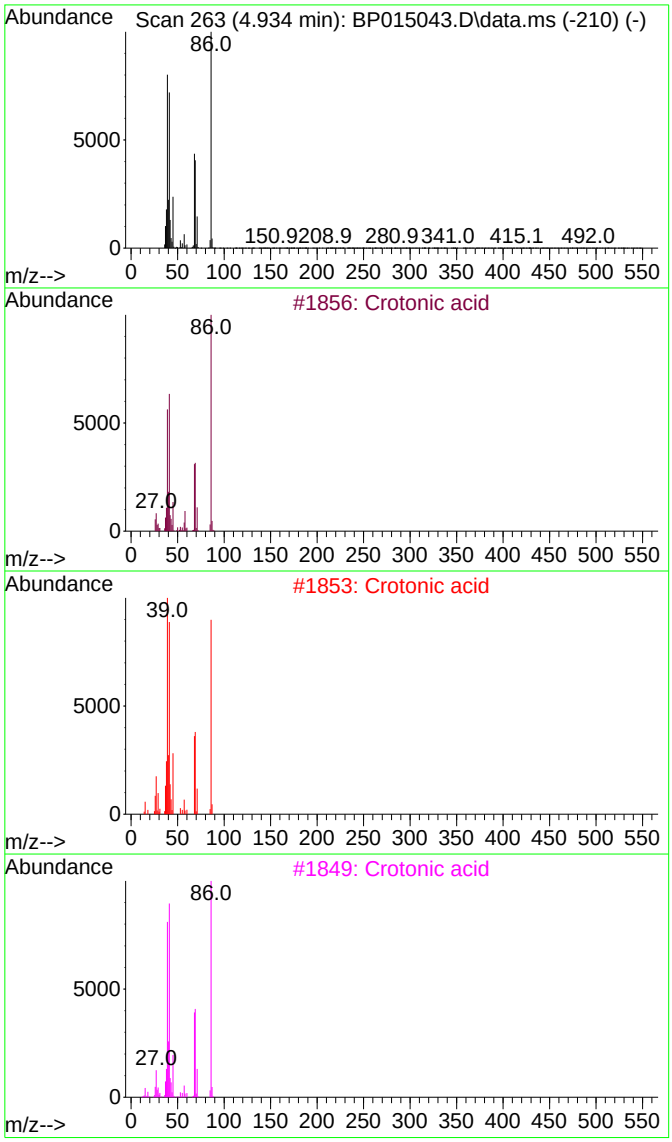
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

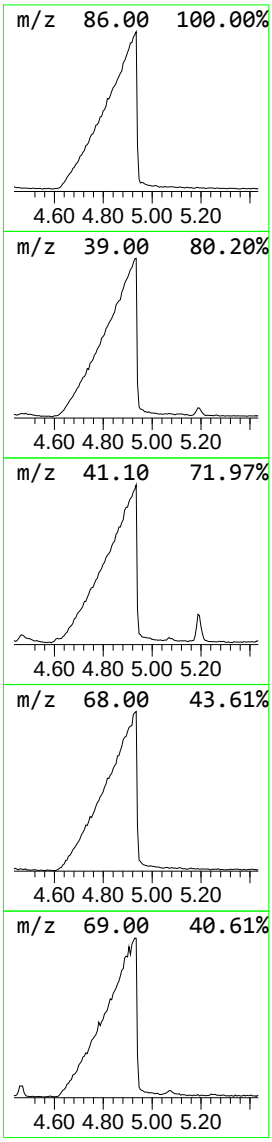
****

Peak Number 6 Crotonic acid Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.934	124.84 ng/ul	6136160	1,4-Dichlorobenzene-d4	8.146

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Crotonic acid	86	C4H6O2	003724-65-0	96
2	Crotonic acid	86	C4H6O2	003724-65-0	96
3	Crotonic acid	86	C4H6O2	003724-65-0	95
4	2-Butenoic acid, (E)-	86	C4H6O2	000107-93-7	94
5	Crotonic acid	86	C4H6O2	003724-65-0	94





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Acq On : 10 May 2023 15:28
Operator : CG/JU
Sample : 02694-18
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
JREP8

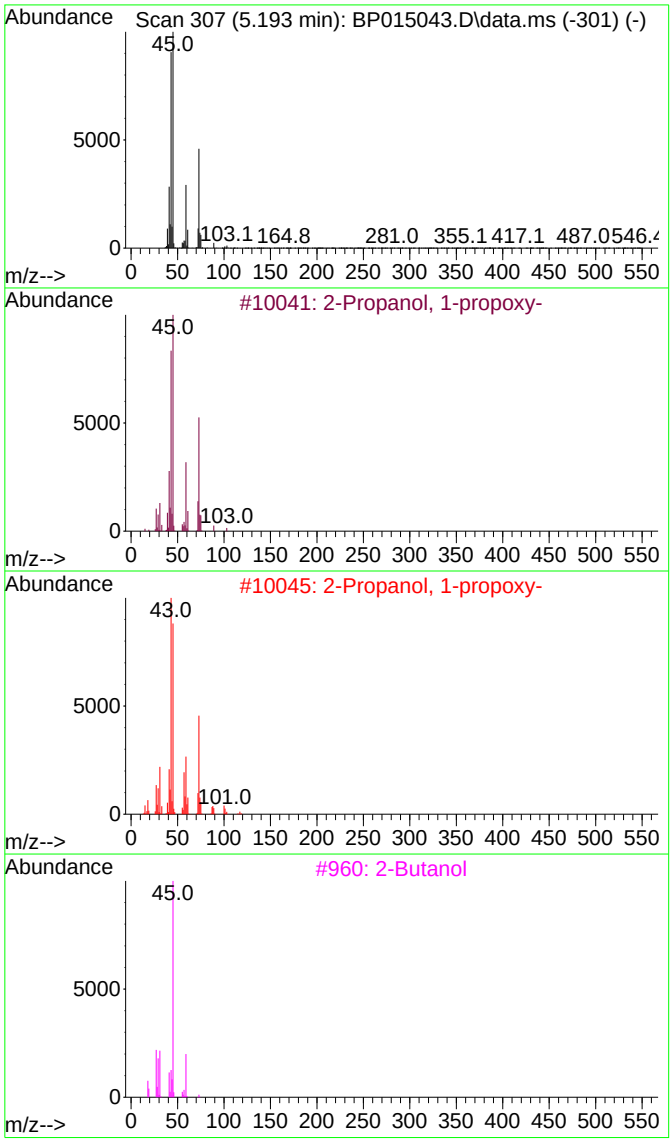
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050323.MA.M
Quant Title : SVOA CALIBRATION

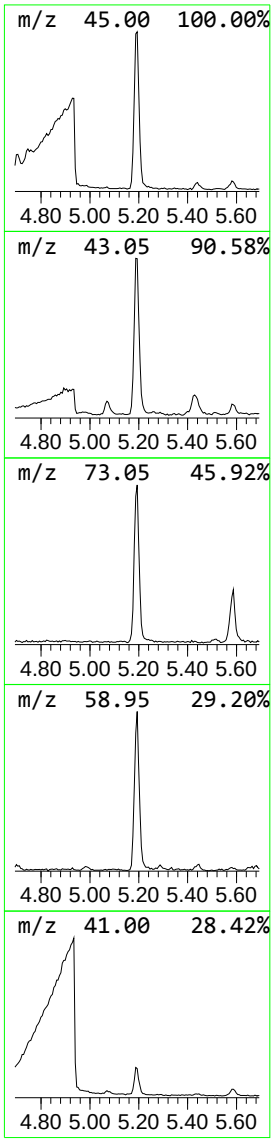
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 2-Propanol, 1-propoxy- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.193	6.78 ng/ul	333058	1,4-Dichlorobenzene-d4	8.146

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Propanol, 1-propoxy-	118	C6H14O2	001569-01-3	86
2	2-Propanol, 1-propoxy-	118	C6H14O2	001569-01-3	64
3	2-Butanol	74	C4H10O	000078-92-2	53
4	2-Propanol, 1-(1-methylethoxy)-	118	C6H14O2	003944-36-3	50
5	2-Propanol, 1-(1-methylethoxy)-	118	C6H14O2	003944-36-3	50





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Sample : 02694-18
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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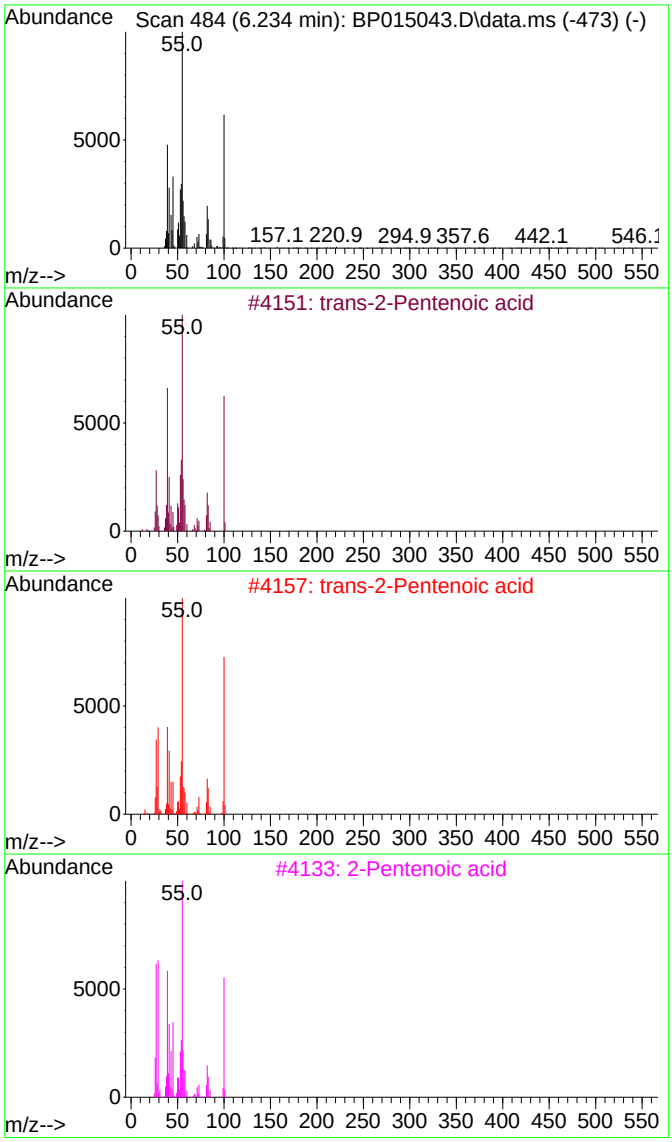
Quant Title : SVOA CALIBRATION

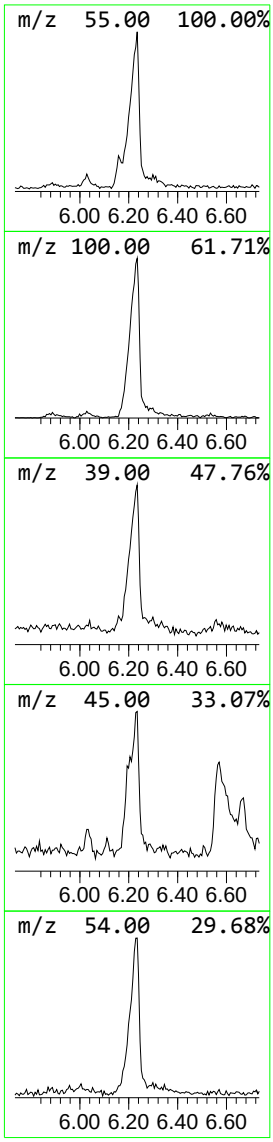
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 trans-2-Pentenoic acid Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.234	5.43 ng/ul	266698	1,4-Dichlorobenzene-d4	8.146

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	trans-2-Pentenoic acid	100	C5H8O2	013991-37-2	95
2	trans-2-Pentenoic acid	100	C5H8O2	013991-37-2	95
3	2-Pentenoic acid	100	C5H8O2	000626-98-2	94
4	Tiglic acid	100	C5H8O2	000080-59-1	87
5	2-Pentenoic acid	100	C5H8O2	000626-98-2	81





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BNA_P
ClientSampleId :
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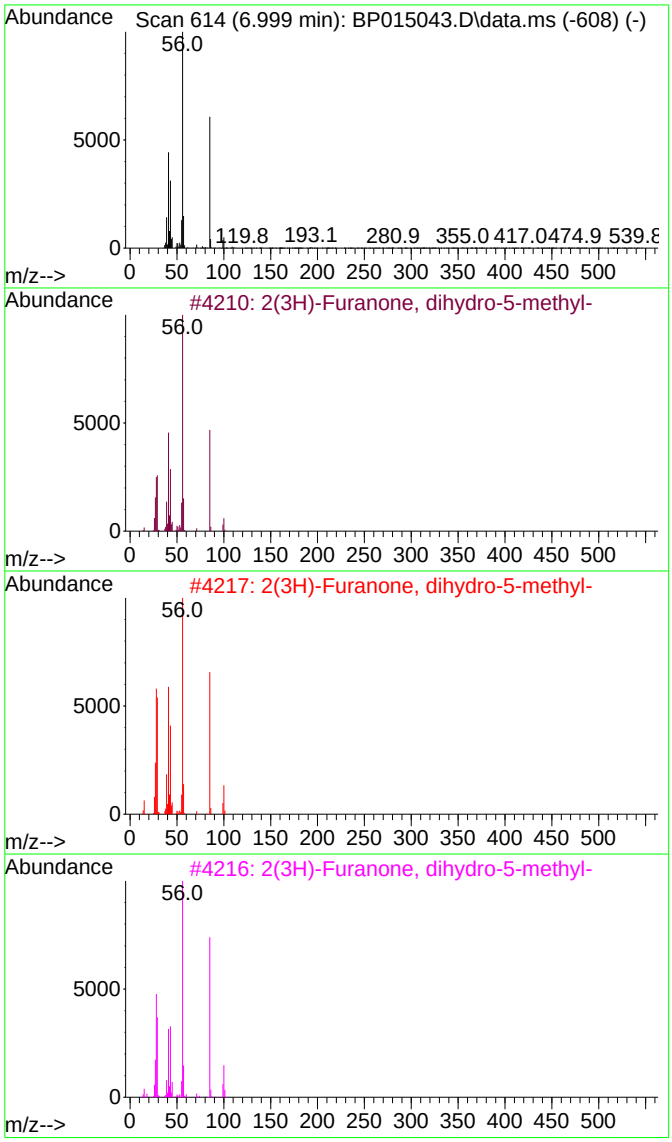
Quant Title : SVOA CALIBRATION

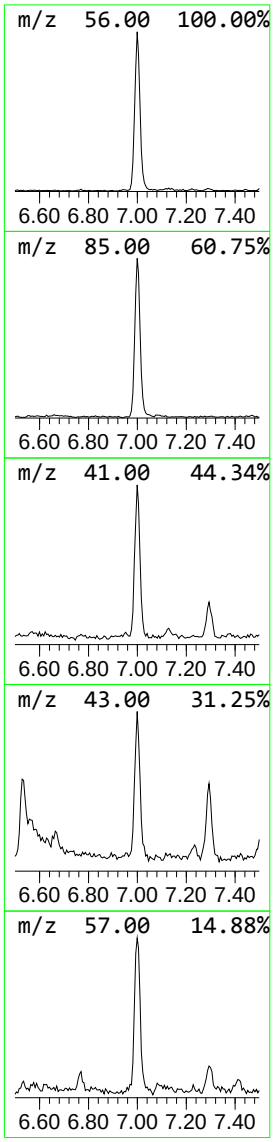
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 2(3H)-Furanone, dihydro-5-m... Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.999	4.25 ng/ul	208730	1,4-Dichlorobenzene-d4	8.146

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2(3H)-Furanone, dihydro-5-methyl-	100	C5H8O2	000108-29-2	91
2	2(3H)-Furanone, dihydro-5-methyl-	100	C5H8O2	000108-29-2	90
3	2(3H)-Furanone, dihydro-5-methyl-	100	C5H8O2	000108-29-2	83
4	2(3H)-Furanone, dihydro-5-methyl-	100	C5H8O2	000108-29-2	72
5	Furan, tetrahydro-2,5-dimethyl-	100	C6H12O	001003-38-9	72





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Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
JREP8

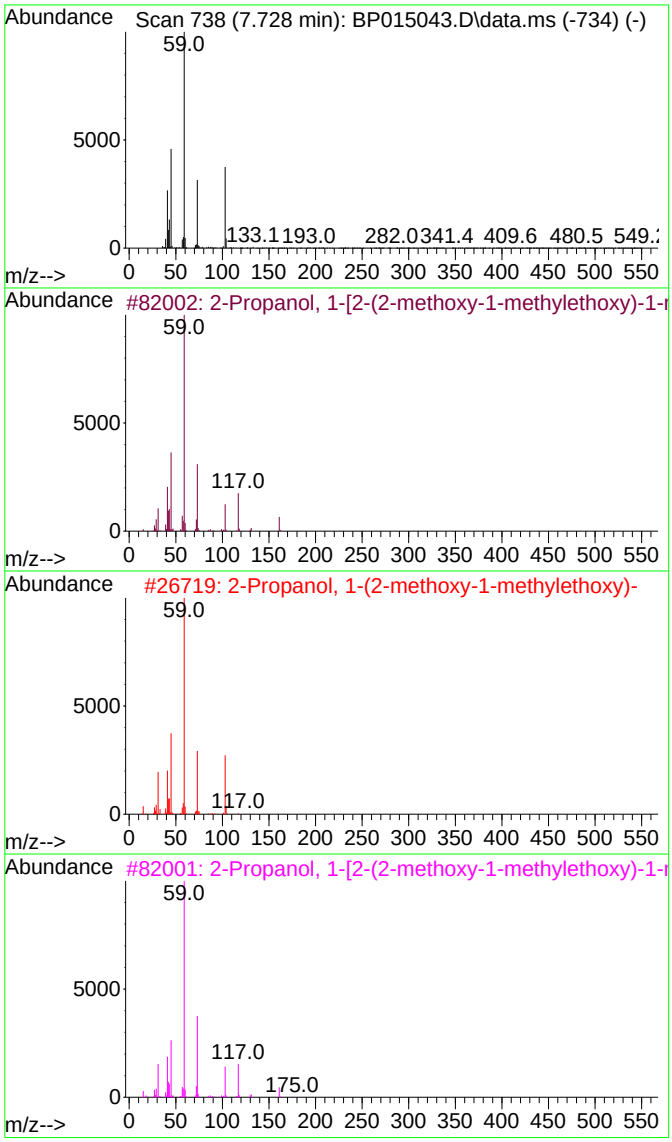
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050323.MA.M
Quant Title : SVOA CALIBRATION

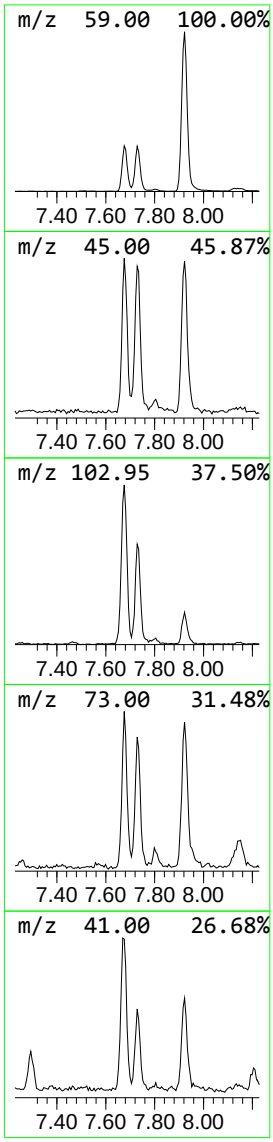
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 2-Propanol, 1-[2-(2-methoxy... Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.728	3.22 ng/ul	158110	1,4-Dichlorobenzene-d4	8.146

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Propanol, 1-[2-(2-methoxy-1-me...	206	C10H22O4	020324-33-8	72
2	2-Propanol, 1-(2-methoxy-1-methy...	148	C7H16O3	020324-32-7	72
3	2-Propanol, 1-[2-(2-methoxy-1-me...	206	C10H22O4	020324-33-8	64
4	1-Propanol, 2-(2-hydroxypropoxy)-	134	C6H14O3	000106-62-7	59
5	Ethyl 2,5,8,11,14-pentaoxahexade...	294	C13H26O7	1000366-80-7	56





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ALS Vial : 4 Sample Multiplier: 1

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BNA_P
ClientSampleId :
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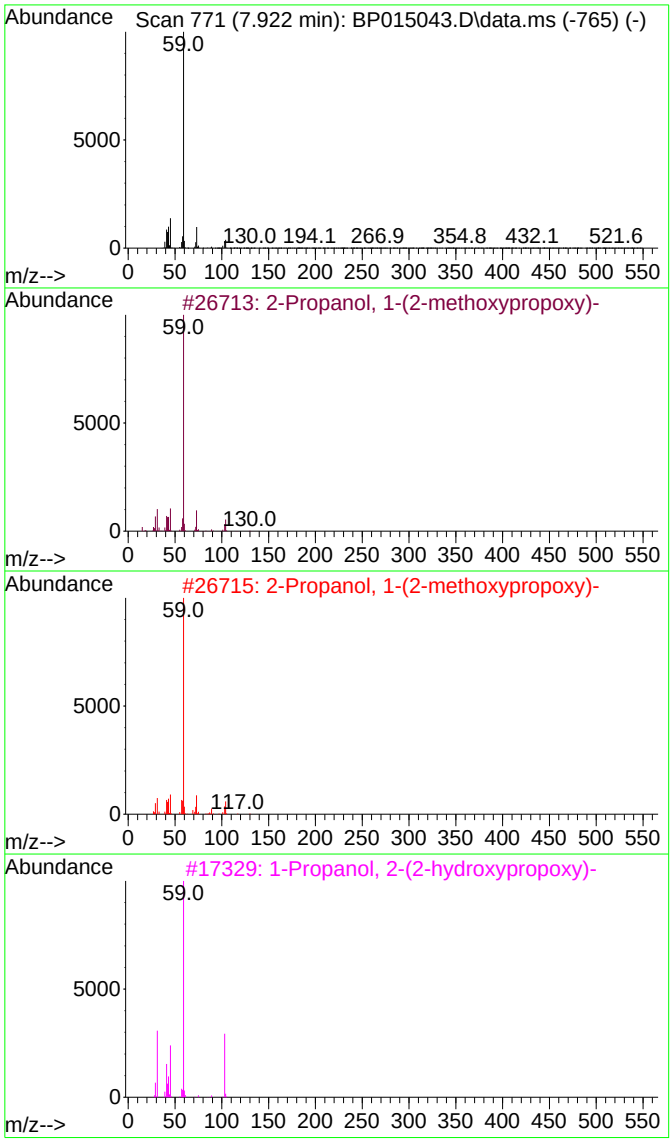
Quant Title : SVOA CALIBRATION

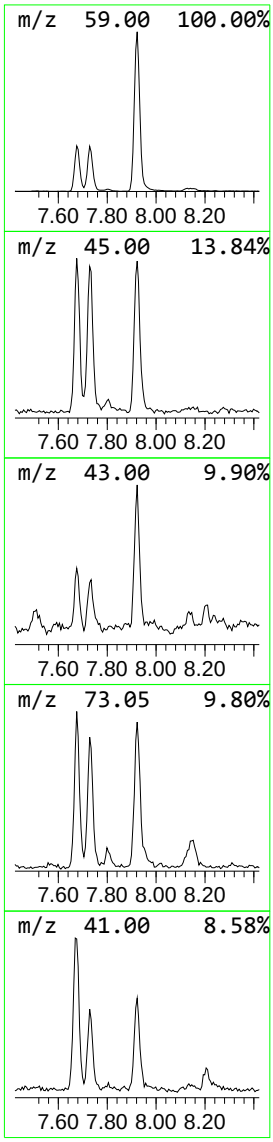
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 11 2-Propanol, 1-(2-methoxypro... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.922	7.04 ng/ul	346024	1,4-Dichlorobenzene-d4	8.146

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Propanol, 1-(2-methoxypropoxy)-	148	C7H16O3	013429-07-7	78
2	2-Propanol, 1-(2-methoxypropoxy)-	148	C7H16O3	013429-07-7	78
3	1-Propanol, 2-(2-hydroxypropoxy)-	134	C6H14O3	000106-62-7	64
4	1-Propanol, 2-(2-hydroxypropoxy)-	134	C6H14O3	000106-62-7	50
5	2-Propanol, 1-(2-methoxy-1-methy...	148	C7H16O3	020324-32-7	50





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Sample : 02694-18
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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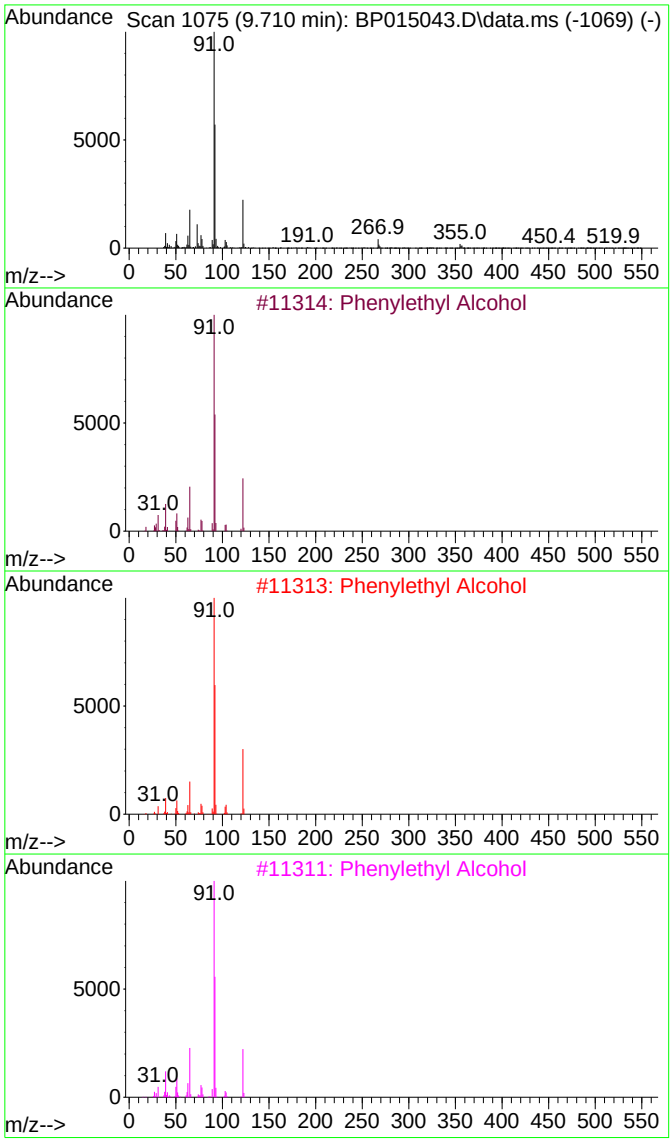
Quant Title : SVOA CALIBRATION

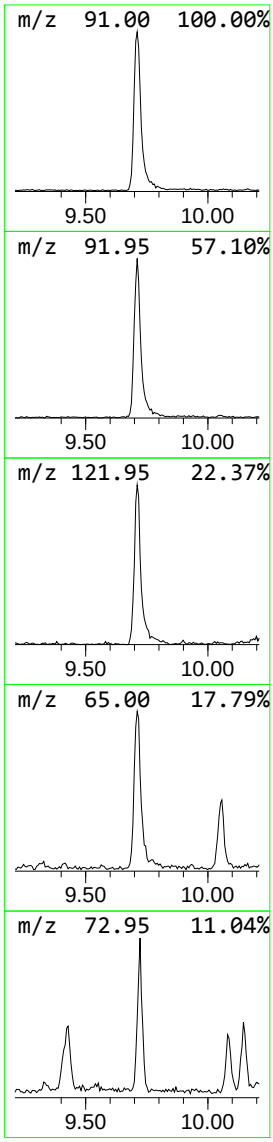
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 12 Phenylethyl Alcohol Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.710	3.28 ng/ul	225602	Naphthalene-d8	10.981

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenylethyl Alcohol	122	C8H10O	000060-12-8	94
2	Phenylethyl Alcohol	122	C8H10O	000060-12-8	93
3	Phenylethyl Alcohol	122	C8H10O	000060-12-8	89
4	Phenylethyl Alcohol	122	C8H10O	000060-12-8	76
5	Spiro[2,4]hepta-4,6-diene	92	C7H8	000765-46-8	49





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Operator : CG/JU
Sample : 02694-18
Misc :
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Instrument :
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ClientSampleId :
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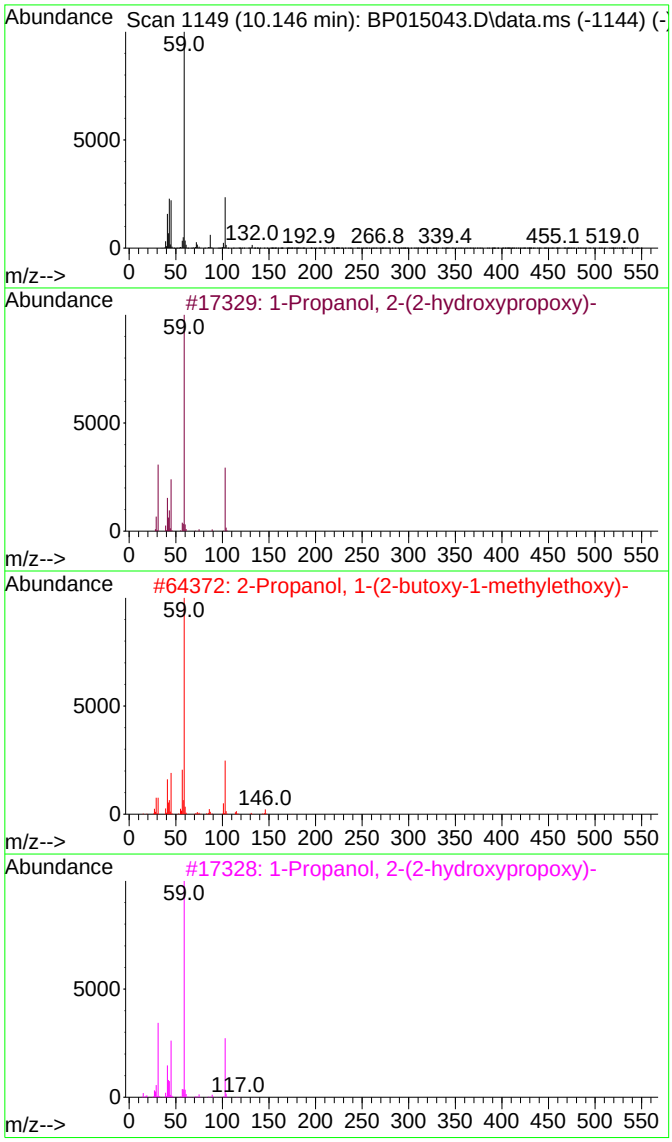
Quant Title : SVOA CALIBRATION

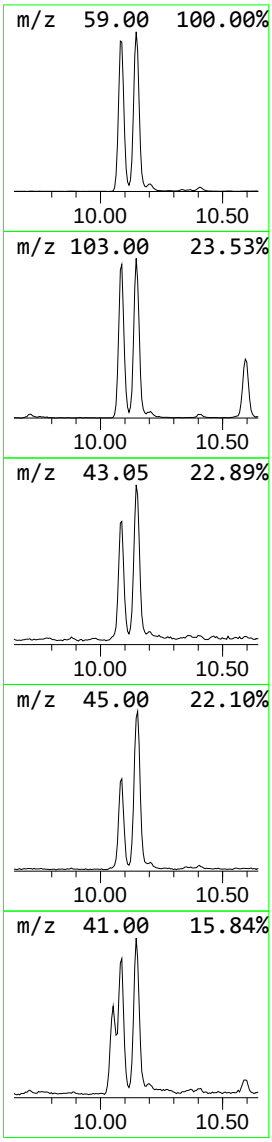
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 13 1-Propanol, 2-(2-hydroxypro... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.146	12.25 ng/ul	843343	Naphthalene-d8	10.981

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Propanol, 2-(2-hydroxypropoxy)-	134	C6H14O3	000106-62-7	78
2	2-Propanol, 1-(2-butoxy-1-methyl...	190	C10H22O3	029911-28-2	78
3	1-Propanol, 2-(2-hydroxypropoxy)-	134	C6H14O3	000106-62-7	78
4	1,4,7-Trimethyl-3,6-dioxaoctane-...	192	C9H20O4	1000423-63-2	72
5	Tetraglyme	222	C10H22O5	000143-24-8	72





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Sample : 02694-18
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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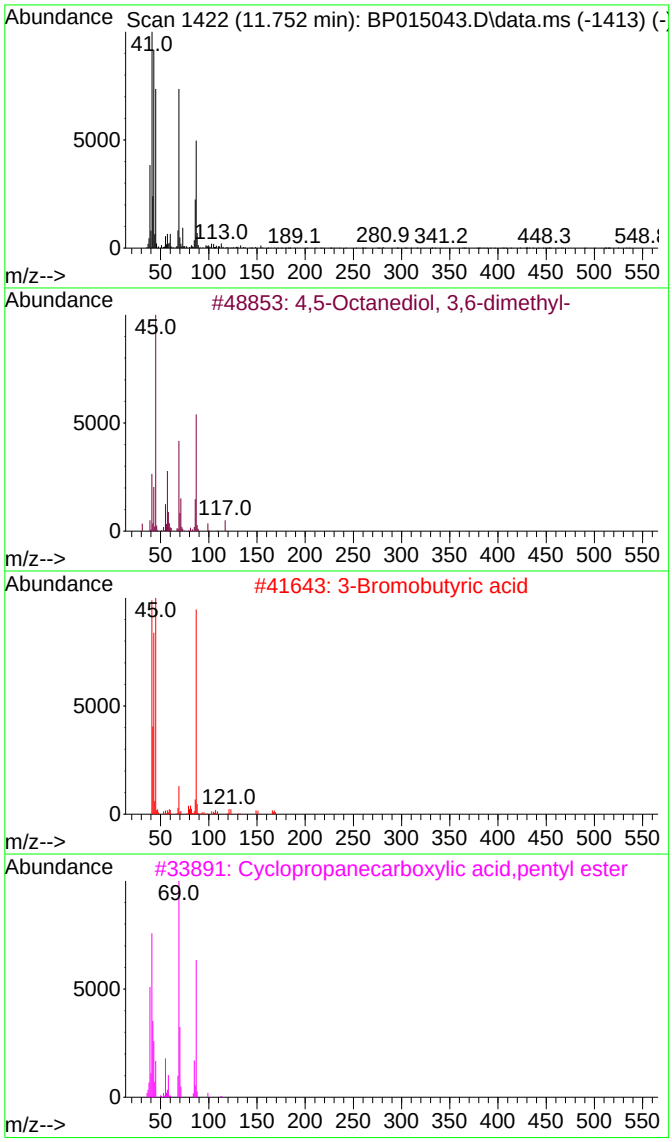
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050323.MA.M
Quant Title : SVOA CALIBRATION

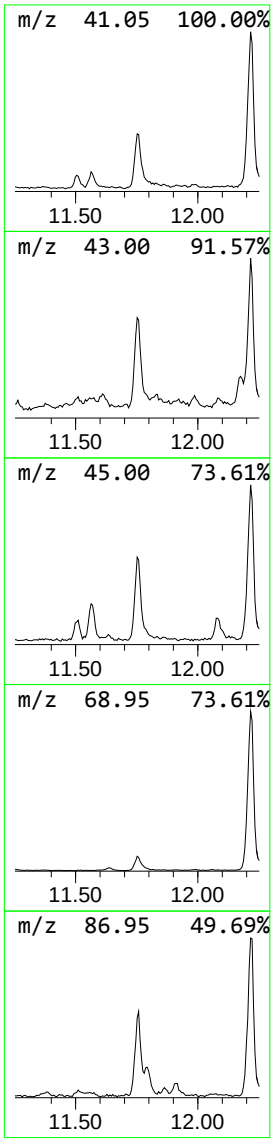
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 14 4,5-Octanediol, 3,6-dimethyl- Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.752	3.60 ng/ul	247451	Naphthalene-d8	10.981

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4,5-Octanediol, 3,6-dimethyl-	174	C10H22O2	1000153-21-1	53
2	3-Bromobutyric acid	166	C4H7BrO2	002623-86-1	43
3	Cyclopropanecarboxylic acid,pent...	156	C9H16O2	060128-02-1	38
4	Cyclopropanecarboxylic acid,isop...	128	C7H12O2	006887-83-8	35
5	2-Propenoic acid, 2-methyl-, dec...	226	C14H26O2	003179-47-3	32





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ALS Vial : 4 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050323.MA.M
Quant Title : SVOA CALIBRATION

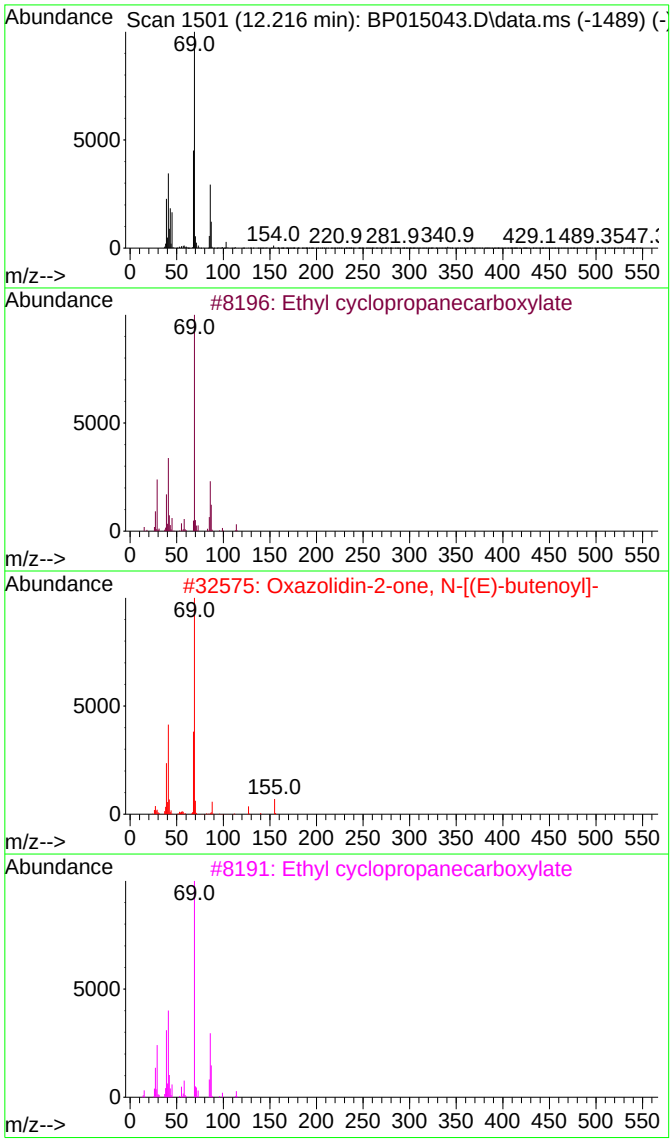
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

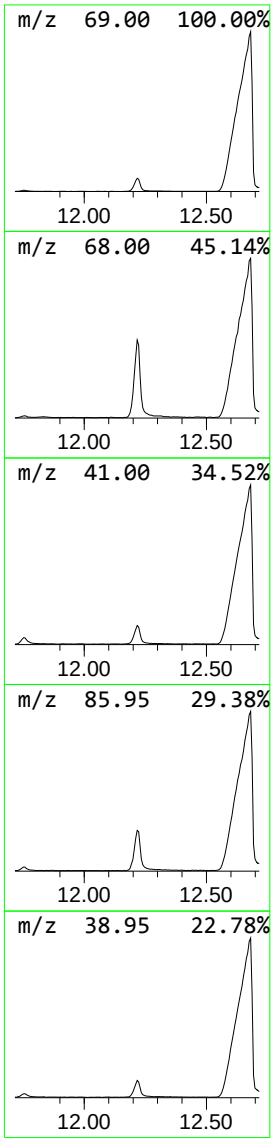
****

Peak Number 15 unknown-02 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.216	13.30 ng/ul	915483	Naphthalene-d8	10.981

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethyl cyclopropanecarboxylate	114	C6H10O2	004606-07-9	42
2	Oxazolidin-2-one, N-[(E)-butenoyl]-	155	C7H9NO3	1010156-45-5	40
3	Ethyl cyclopropanecarboxylate	114	C6H10O2	004606-07-9	40
4	Cyclopropanecarboxylic acid, 3-m...	154	C9H14O2	1000299-37-4	36
5	1,5-Pentanediol, O,O'-di(proparg...	268	C13H16O6	1000331-35-4	36





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ALS Vial : 4 Sample Multiplier: 1

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ClientSampleId :
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Quant Title : SVOA CALIBRATION

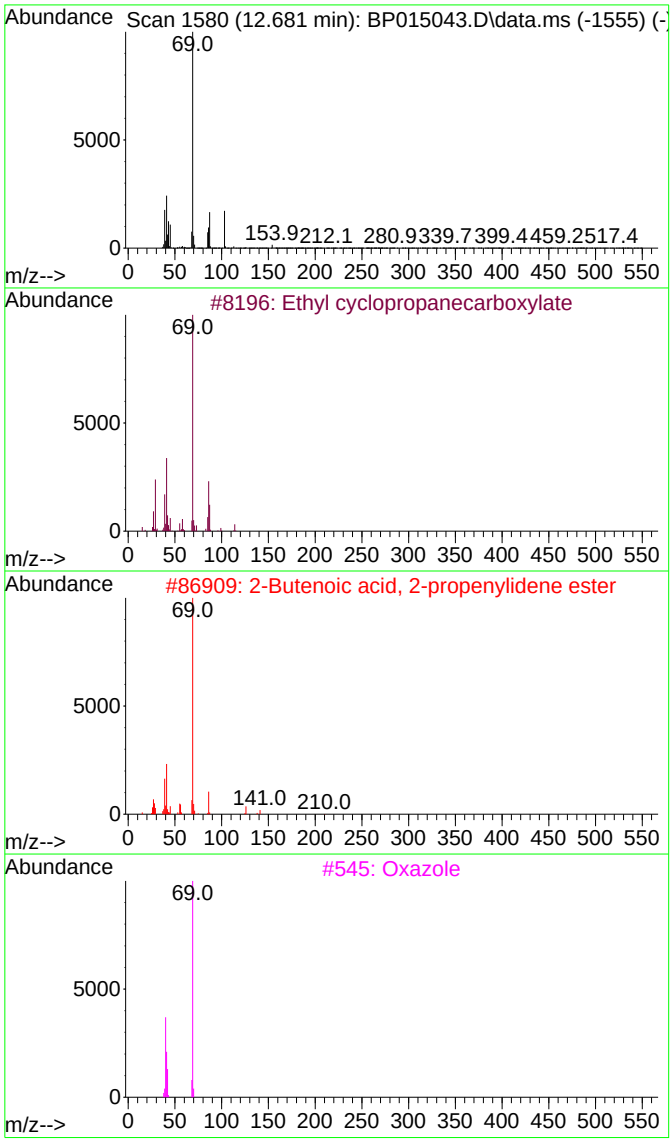
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

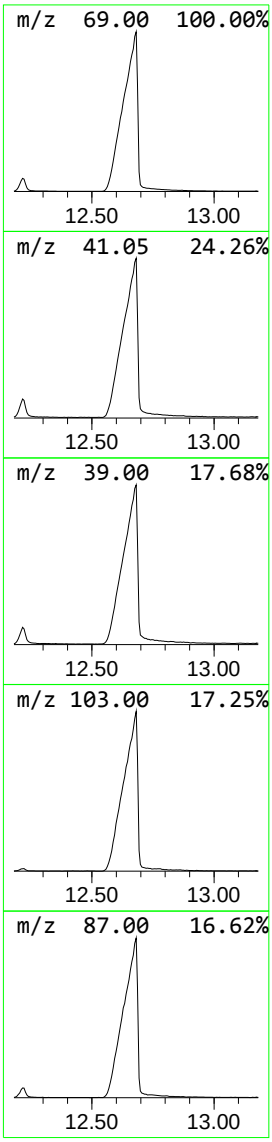
****

Peak Number 16 unknown-03 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.681	285.41 ng/ul	19643500	Naphthalene-d8	10.981

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethyl cyclopropanecarboxylate	114	C6H10O2	004606-07-9	38
2	2-Butenoic acid, 2-propenylidene...	210	C11H14O4	055030-70-1	36
3	Oxazole	69	C3H3NO	000288-42-6	9
4	Cyclopropanecarboxylic acid, 3-m...	154	C9H14O2	1000299-37-4	9
5	Cyclopropanecarboxylic acid,buty...	142	C8H14O2	054947-39-6	9





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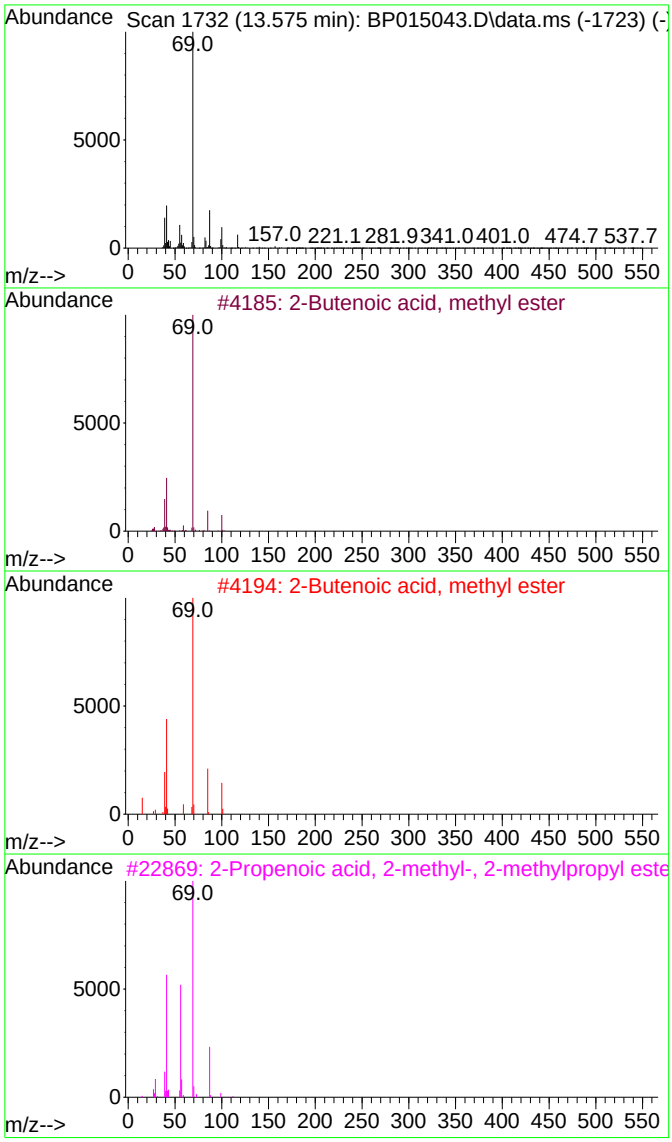
Quant Title : SVOA CALIBRATION

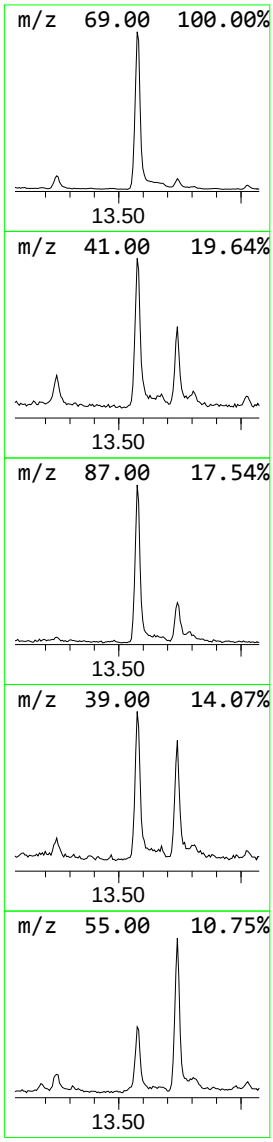
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 17 2-Butenoic acid, methyl ester Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.575	7.50 ng/ul	694314	Acenaphthene-d10	14.787

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Butenoic acid, methyl ester	100	C5H8O2	018707-60-3	59
2	2-Butenoic acid, methyl ester	100	C5H8O2	018707-60-3	45
3	2-Propenoic acid, 2-methyl-, 2-m...	142	C8H14O2	000097-86-9	45
4	4-Hexen-3-one	98	C6H10O	002497-21-4	40
5	2-Hydroxyethyl methacrylate	130	C6H10O3	000868-77-9	40





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Data File : BP015043.D
Acq On : 10 May 2023 15:28
Operator : CG/JU
Sample : 02694-18
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
JREP8

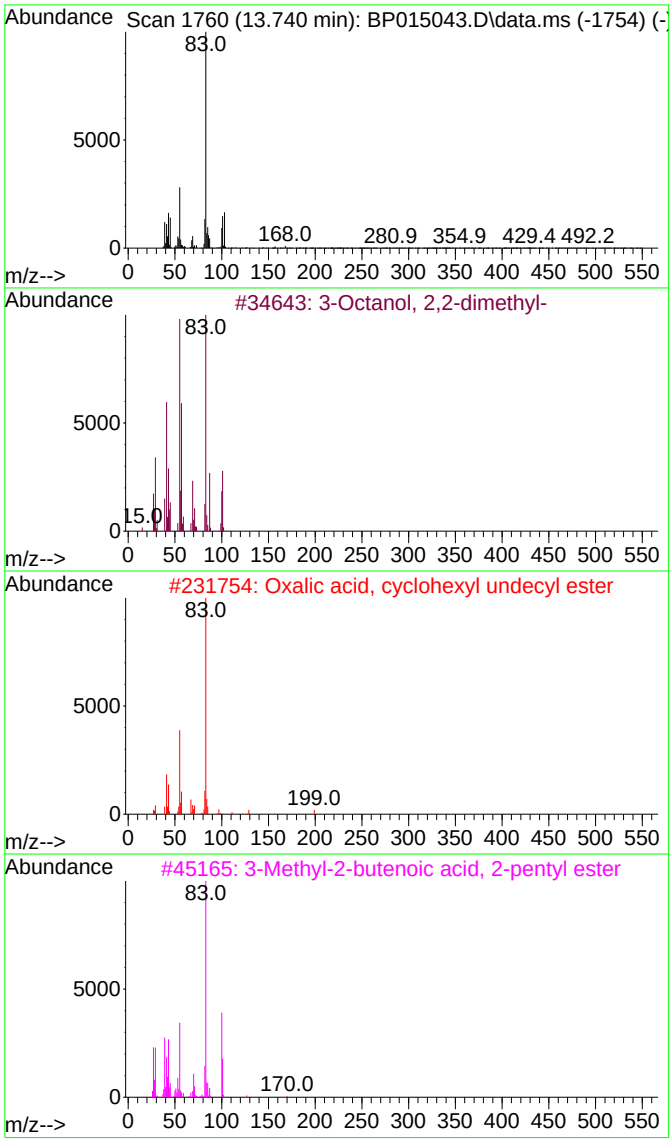
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050323.MA.M
Quant Title : SVOA CALIBRATION

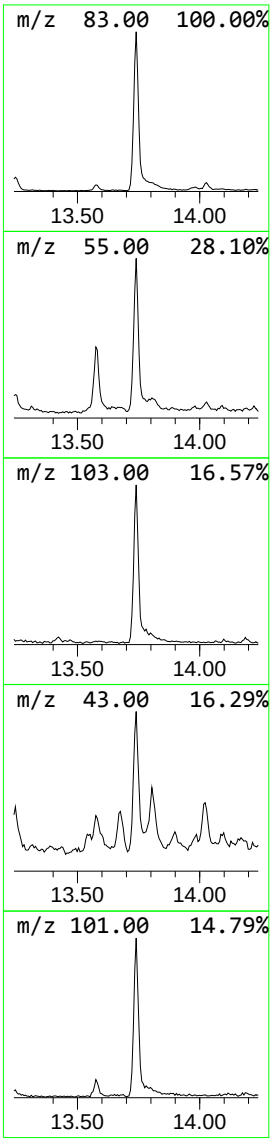
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 18 3-Octanol, 2,2-dimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.740	7.82 ng/ul	723340	Acenaphthene-d10	14.787

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Octanol, 2,2-dimethyl-	158	C10H22O	019841-72-6	59
2	Oxalic acid, cyclohexyl undecyl ...	326	C19H34O4	1000309-31-3	47
3	3-Methyl-2-butenoic acid, 2-pent...	170	C10H18O2	150462-84-3	45
4	2-Butenoic acid, 3-methyl-, ethy...	128	C7H12O2	000638-10-8	45
5	3-Methyl-2-butenoic acid, cyclop...	168	C10H16O2	1000279-23-3	45





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Operator : CG/JU
Sample : 02694-18
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
JREP8

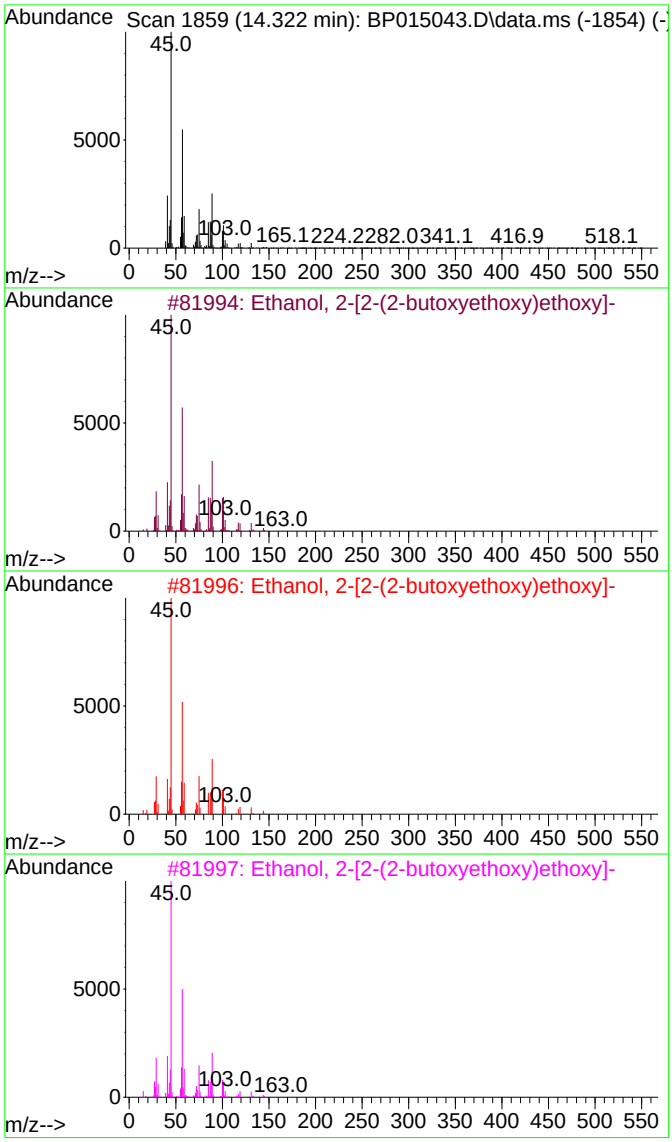
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050323.MA.M
Quant Title : SVOA CALIBRATION

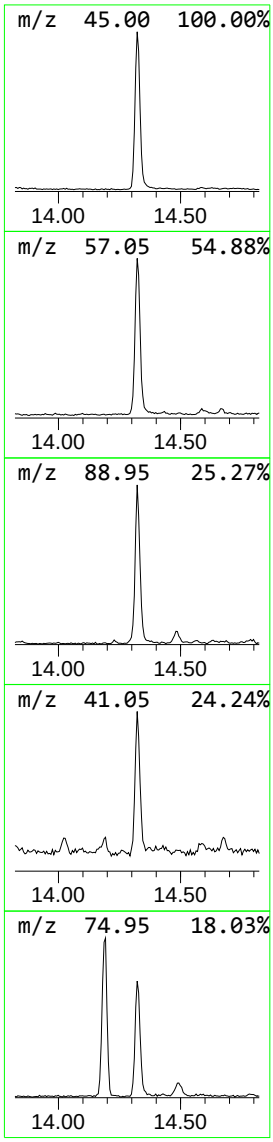
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 19 Ethanol, 2-[2-(2-butoxyetho... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.322	5.66 ng/ul	523731	Acenaphthene-d10	14.787

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethanol, 2-[2-(2-butoxyethoxy)et...	206	C10H22O4	000143-22-6	91
2	Ethanol, 2-[2-(2-butoxyethoxy)et...	206	C10H22O4	000143-22-6	91
3	Ethanol, 2-[2-(2-butoxyethoxy)et...	206	C10H22O4	000143-22-6	87
4	3,6,9,12-Tetraoxahexadecan-1-ol	250	C12H26O5	001559-34-8	86
5	Ethanol, 2-[2-(2-butoxyethoxy)et...	206	C10H22O4	000143-22-6	83





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Sample : 02694-18
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
JREP8

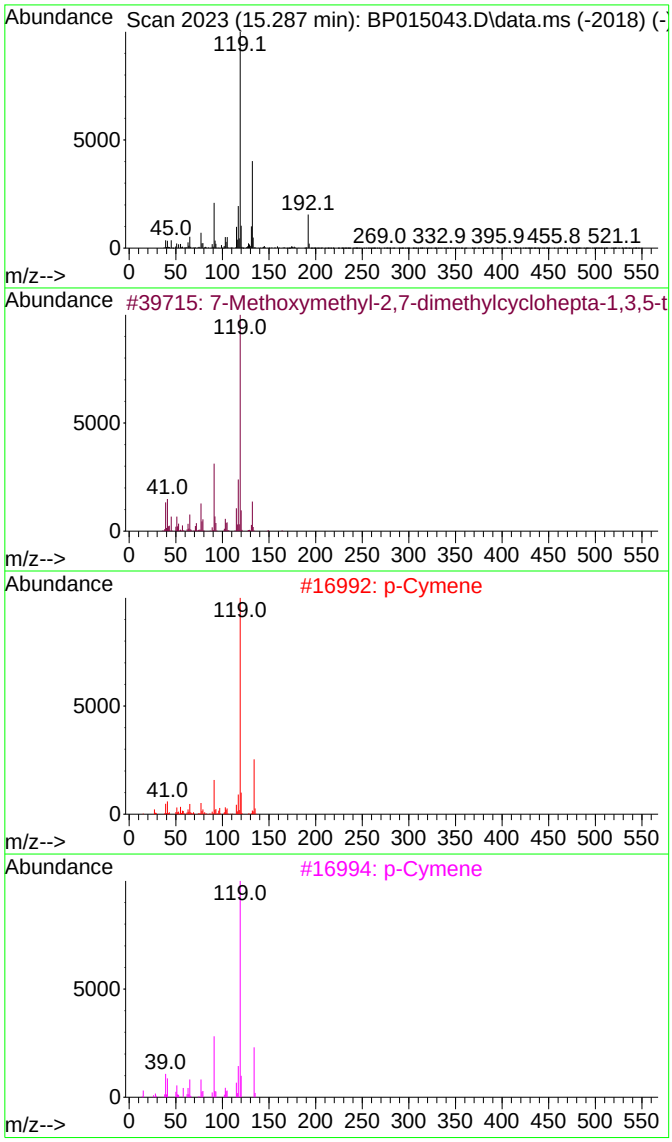
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050323.MA.M
Quant Title : SVOA CALIBRATION

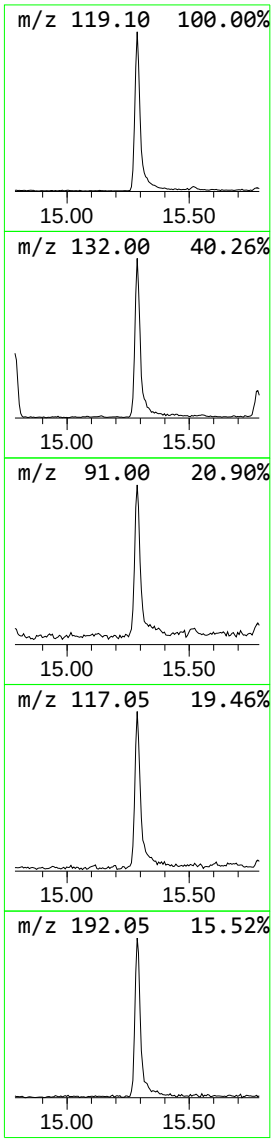
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 20 7-Methoxymethyl-2,7-dimethy... Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.287	4.05 ng/ul	374889	Acenaphthene-d10	14.787

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	7-Methoxymethyl-2,7-dimethylcycl...	164	C11H16O	073992-48-0	59
2	p-Cymene	134	C10H14	000099-87-6	43
3	p-Cymene	134	C10H14	000099-87-6	43
4	p-Cymene	134	C10H14	000099-87-6	43
5	o-Cymene	134	C10H14	000527-84-4	43





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Acq On : 10 May 2023 15:28
Operator : CG/JU
Sample : 02694-18
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
JREP8

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050323.MA.M
Quant Title : SVOA CALIBRATION

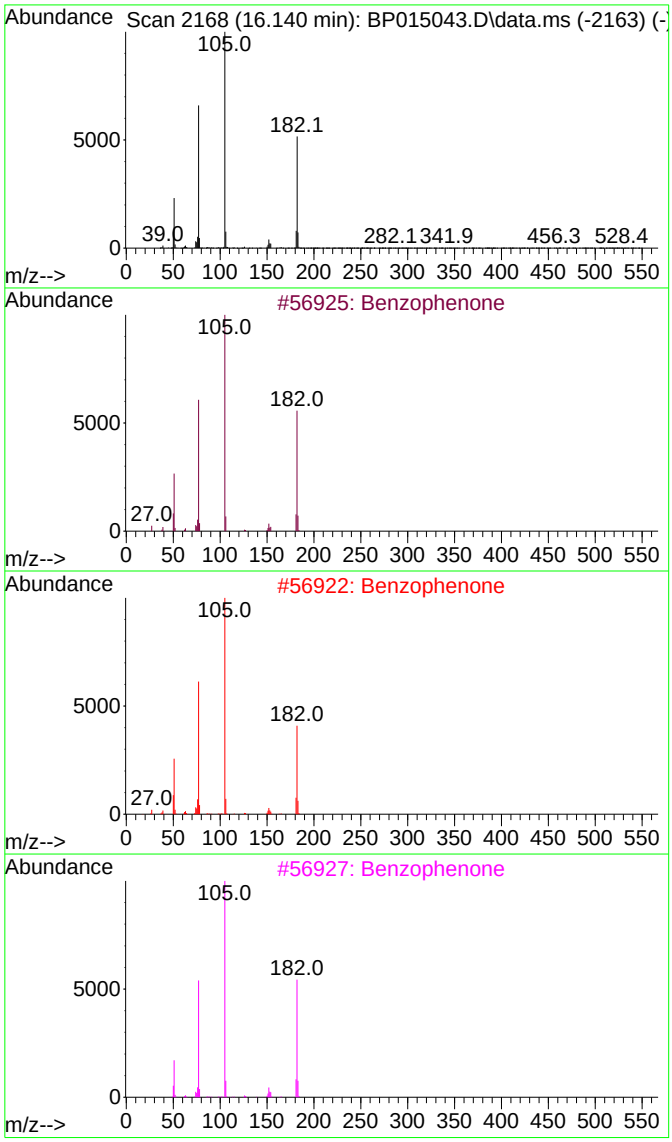
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

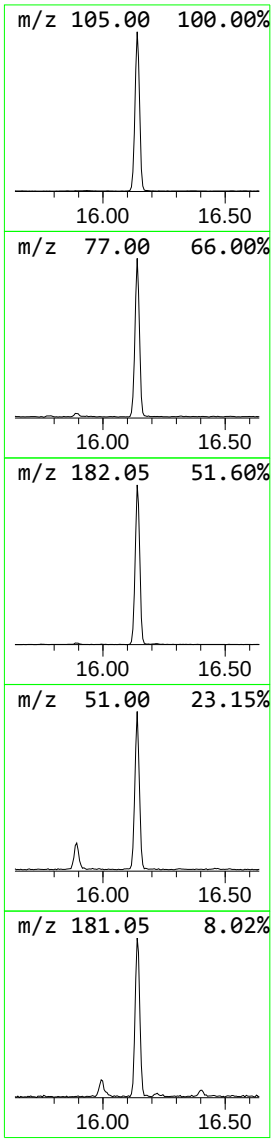
****

Peak Number 21 Benzophenone Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.140	11.90 ng/ul	1101050	Acenaphthene-d10	14.787

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzophenone	182	C13H10O	000119-61-9	96
2	Benzophenone	182	C13H10O	000119-61-9	95
3	Benzophenone	182	C13H10O	000119-61-9	95
4	Benzophenone	182	C13H10O	000119-61-9	93
5	1,2,4,5-Tetroxane, 3,3,6,6-tetra...	396	C26H20O4	016204-36-7	64





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Operator : CG/JU
Sample : 02694-18
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
JREP8

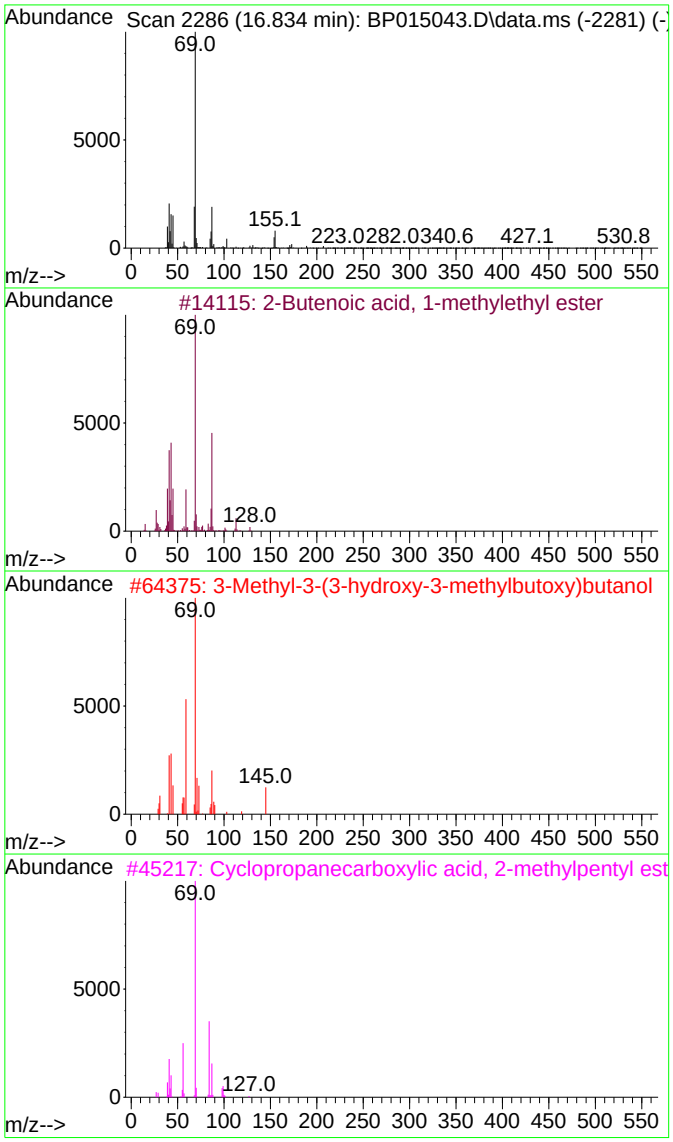
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050323.MA.M
Quant Title : SVOA CALIBRATION

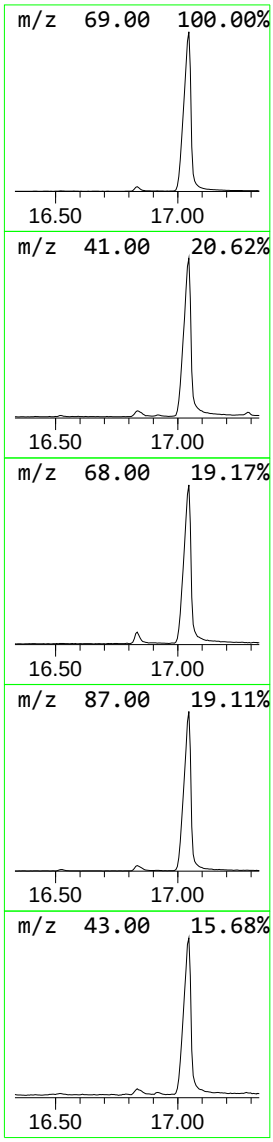
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 22 unknown-04 Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.834	4.49 ng/ul	590975	Phenanthrene-d10	17.540

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Butenoic acid, 1-methylethyl e...	128	C7H12O2	018060-77-0	42
2	3-Methyl-3-(3-hydroxy-3-methylbu...	190	C10H22O3	1000432-16-8	39
3	Cyclopropanecarboxylic acid, 2-m...	170	C10H18O2	1000354-65-8	38
4	Oxazolidin-2-one, N-[(E)-butenoyl]-	155	C7H9NO3	1010156-45-5	38
5	3-Heptanol, 2-methyl-	130	C8H18O	018720-62-2	28





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Acq On : 10 May 2023 15:28
Operator : CG/JU
Sample : 02694-18
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
JREP8

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050323.MA.M
Quant Title : SVOA CALIBRATION

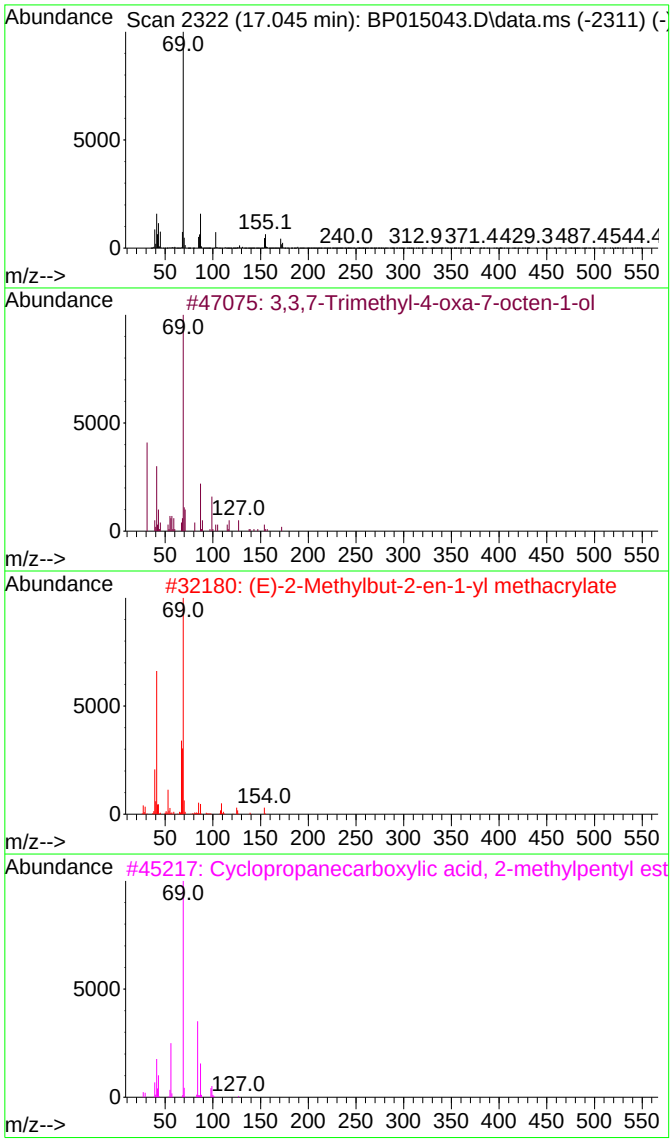
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

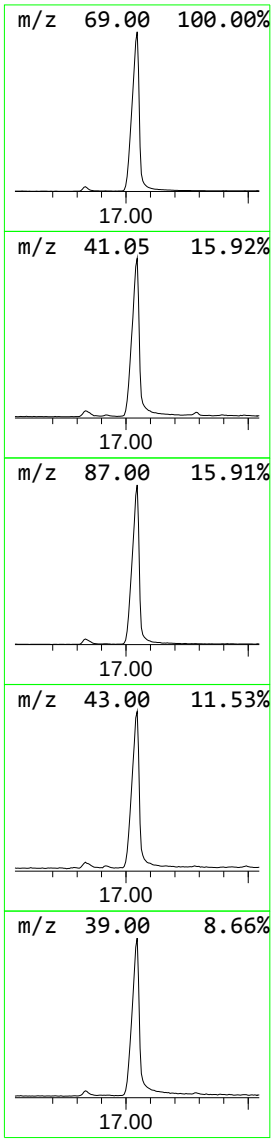
****

Peak Number 23 unknown-05 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.045	115.04 ng/ul	15150700	Phenanthrene-d10	17.540

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3,3,7-Trimethyl-4-oxa-7-octen-1-ol	172	C10H20O2	069161-47-3	39
2	(E)-2-Methylbut-2-en-1-yl methac...	154	C9H14O2	088142-95-4	38
3	Cyclopropanecarboxylic acid, 2-m...	170	C10H18O2	1000354-65-8	36
4	Ethyl cyclopropanecarboxylate	114	C6H10O2	004606-07-9	36
5	Oxazolidin-2-one, N-[(E)-butenoyl]-	155	C7H9NO3	1010156-45-5	33





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Operator : CG/JU
Sample : 02694-18
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

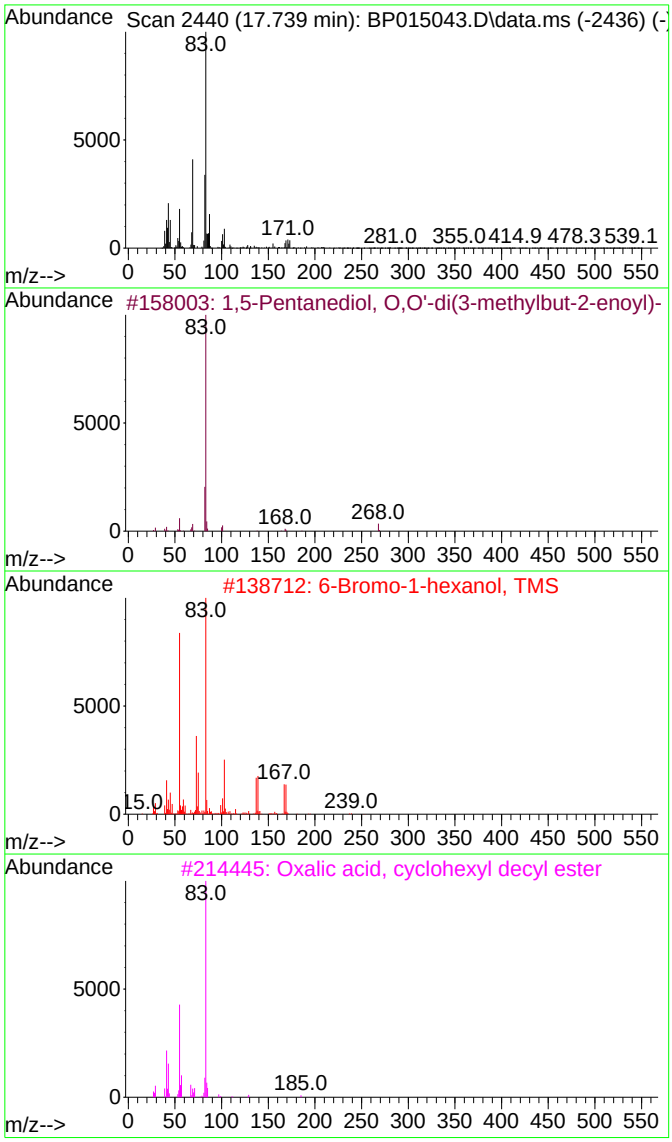
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

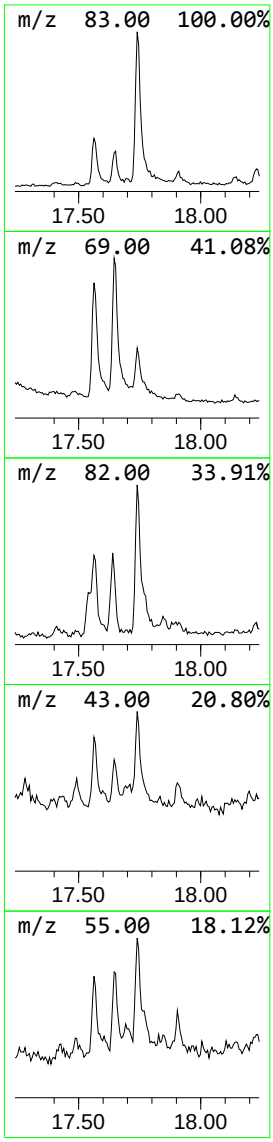
****

Peak Number 24 unknown-06 Concentration Rank 29

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.740	2.63 ng/ul	345736	Phenanthrene-d10	17.540

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,5-Pentanediol, O,O'-di(3-methy...	268	C15H24O4	1010331-15-6	42
2	6-Bromo-1-hexanol, TMS	252	C9H21BrOSi	026306-00-3	38
3	Oxalic acid, cyclohexyl decyl ester	312	C18H32O4	1000309-31-2	38
4	Oxalic acid, cyclohexyl isohexyl...	256	C14H24O4	1010309-30-7	38
5	2-Butenoic acid, 3-methyl-, 3-me...	168	C10H16O2	072779-06-7	38





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Sample : 02694-18
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
JREP8

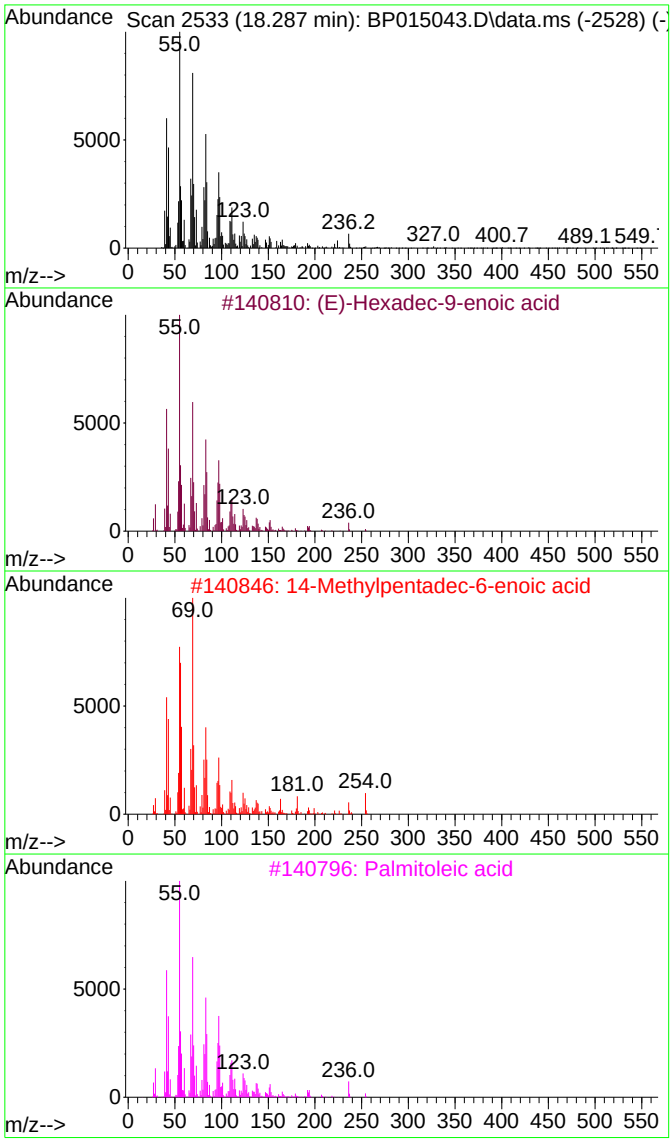
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050323.MA.M
Quant Title : SVOA CALIBRATION

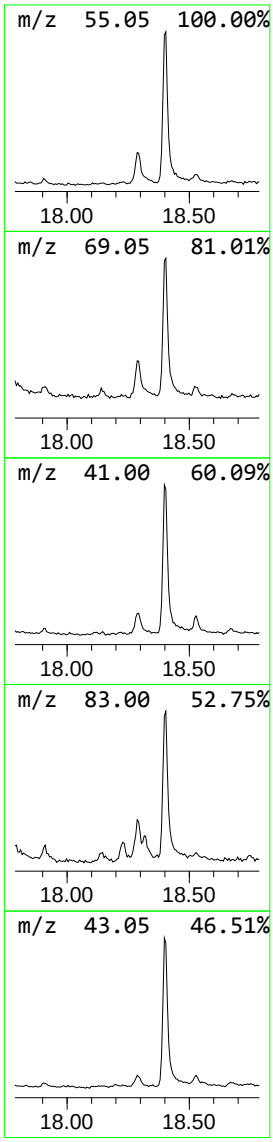
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 25 (E)-Hexadec-9-enoic acid Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.287	2.83 ng/ul	373146	Phenanthrene-d10	17.540

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	(E)-Hexadec-9-enoic acid	254	C16H30O2	010030-73-6	97
2	14-Methylpentadec-6-enoic acid	254	C16H30O2	127486-79-7	95
3	Palmitoleic acid	254	C16H30O2	000373-49-9	90
4	Erucic acid	338	C22H42O2	000112-86-7	90
5	2-Undecanol oleate	436	C29H56O2	1000465-66-9	87





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Misc :
ALS Vial : 4 Sample Multiplier: 1

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ClientSampleId :
JREP8

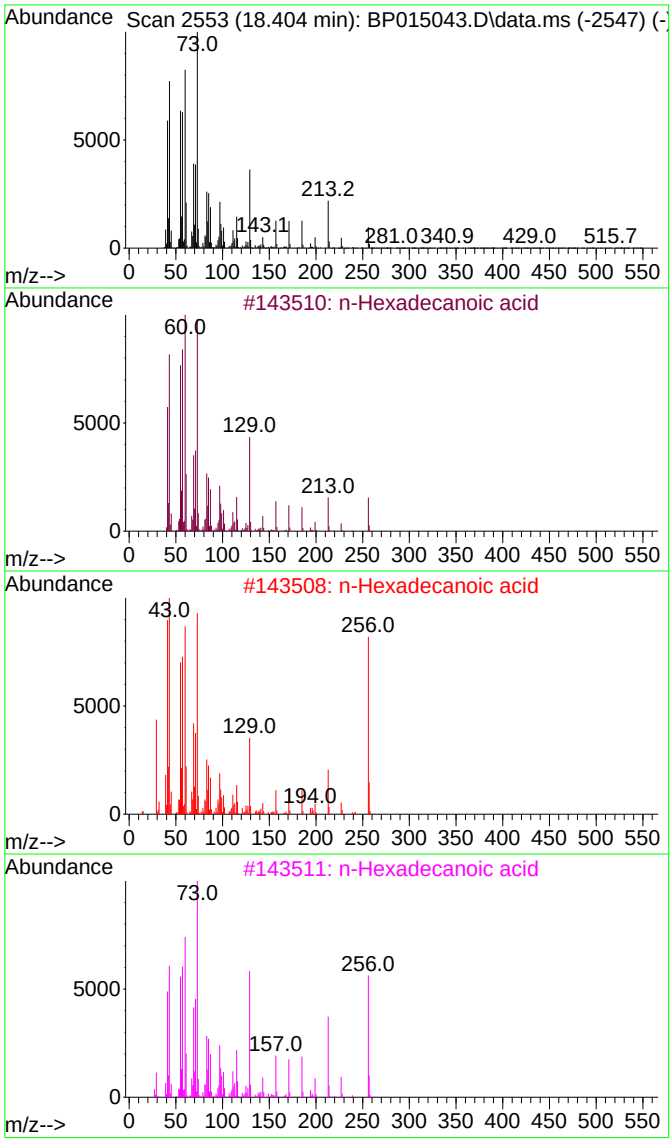
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050323.MA.M
Quant Title : SVOA CALIBRATION

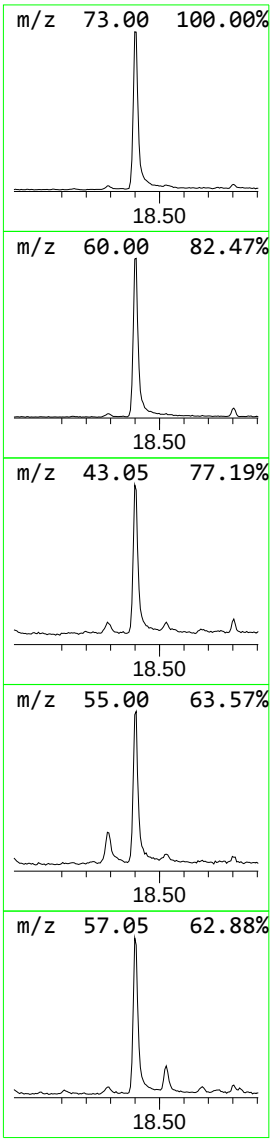
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 26 n-Hexadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.404	16.94 ng/ul	2231490	Phenanthrene-d10	17.540

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
3	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
4	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	96





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Misc :
ALS Vial : 4 Sample Multiplier: 1

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ClientSampleId :
JREP8

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Quant Title : SVOA CALIBRATION

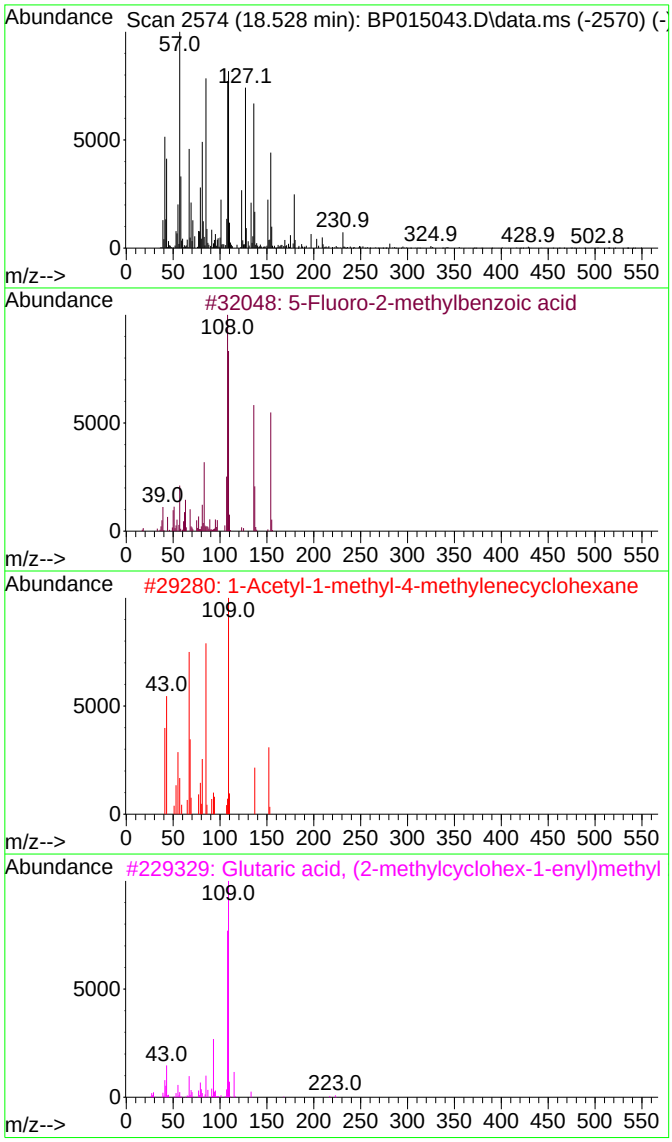
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

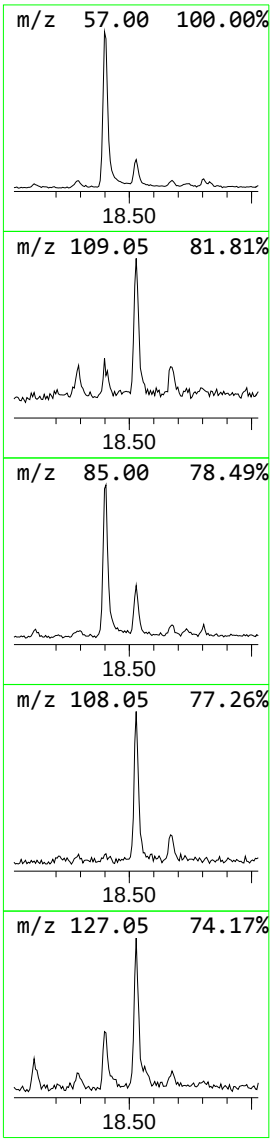
****

Peak Number 27 unknown-07 Concentration Rank 28

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.528	2.64 ng/ul	347176	Phenanthrene-d10	17.540

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	5-Fluoro-2-methylbenzoic acid	154	C8H7FO2	033184-16-6	45
2	1-Acetyl-1-methyl-4-methylenecyc...	152	C10H16O	1000424-28-4	22
3	Glutaric acid, (2-methylcyclohex...	324	C19H32O4	1000405-50-3	22
4	Benzoic acid, 2,6-dihydroxy-	154	C7H6O4	000303-07-1	22
5	Glutaric acid, (2-methylcyclohex...	334	C19H23FO4	1000405-50-5	14





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ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

****
Peak Number 28 Octadecanoic acid Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.622	9.44 ng/ul	1302130	Chrysene-d12	21.622

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecanoic acid	284	C18H36O2	000057-11-4	99
2	Octadecanoic acid	284	C18H36O2	000057-11-4	99
3	Octadecanoic acid	284	C18H36O2	000057-11-4	99
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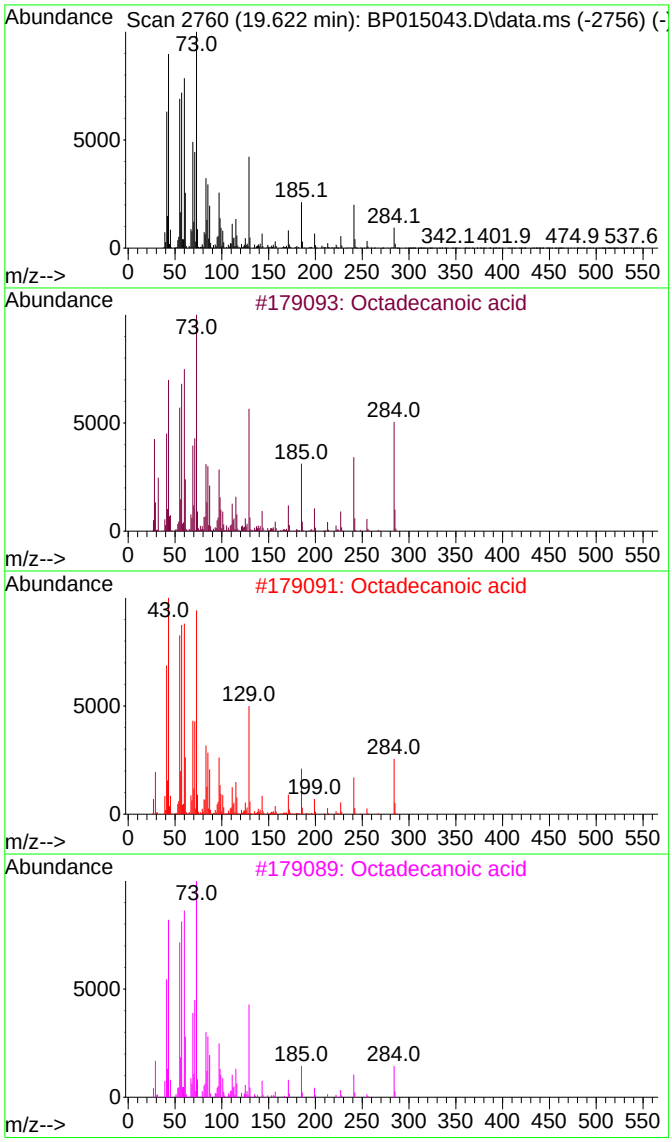
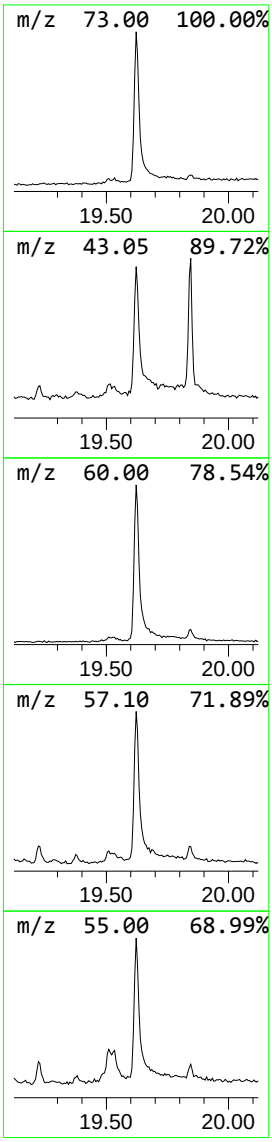
5	Pentadecanoic acid	242	C15H30O2	001002-84-2	90





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Acq On : 10 May 2023 15:28
Operator : CG/JU
Sample : 02694-18
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
JREP8

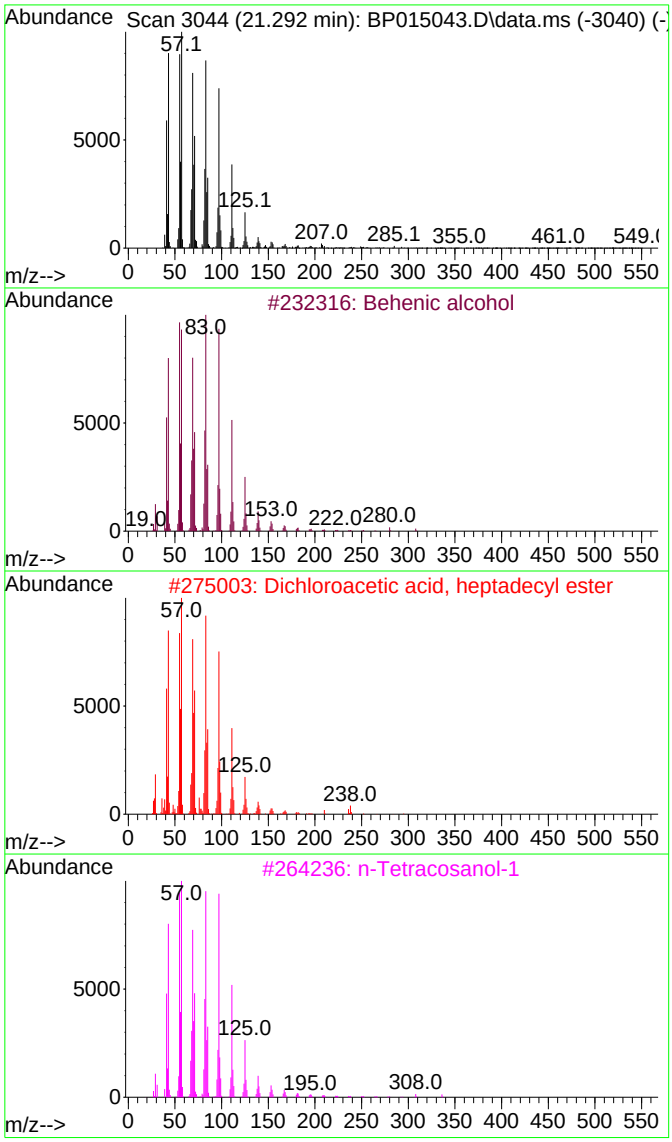
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050323.MA.M
Quant Title : SVOA CALIBRATION

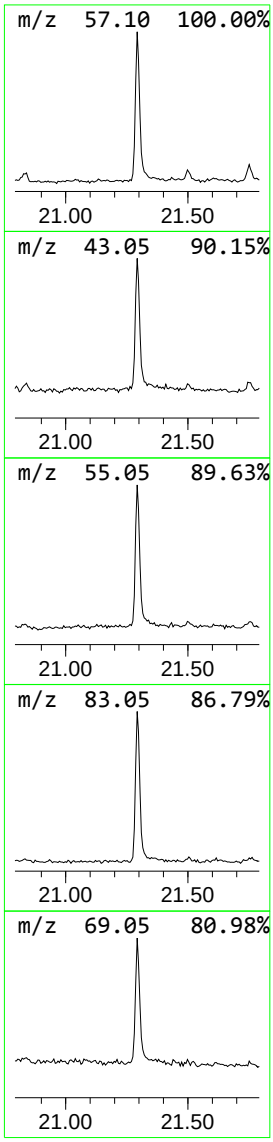
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 29 Behenic alcohol Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.292	6.36 ng/ul	877517	Chrysene-d12	21.622

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Behenic alcohol	326	C22H46O	000661-19-8	94
2	Dichloroacetic acid, heptadecyl ...	366	C19H36Cl2O2	1000282-98-2	94
3	n-Tetracosanol-1	354	C24H50O	000506-51-4	94
4	5-Octadecene, (E)-	252	C18H36	007206-21-5	93
5	Cyclopentadecane	210	C15H30	000295-48-7	93





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051023\
Data File : BP015043.D
Acq On : 10 May 2023 15:28
Operator : CG/JU
Sample : 02694-18
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
JREP8

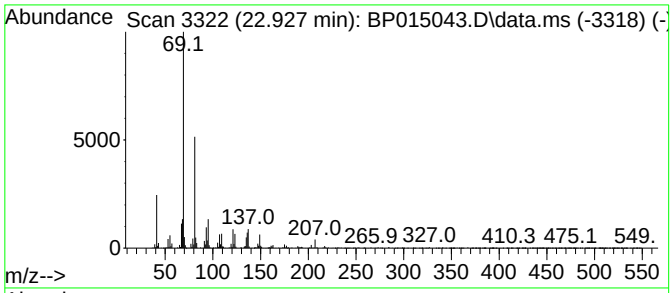
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050323.MA.M
Quant Title : SVOA CALIBRATION

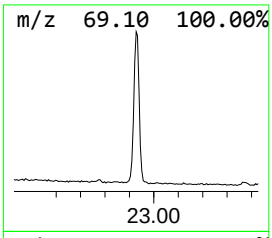
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

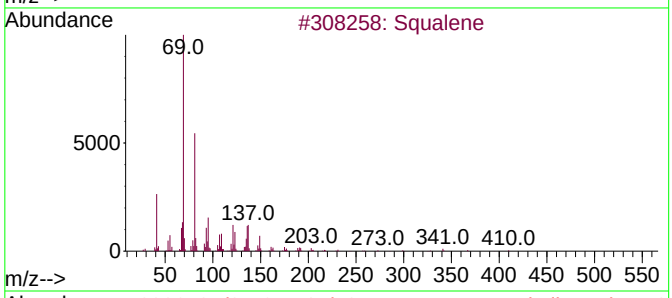
****
Peak Number 30 Squalene Concentration Rank 19

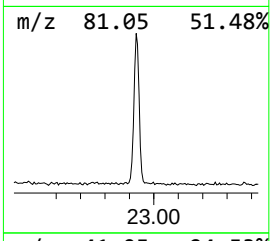
R.T.	EstConc	Area	Relative to ISTD	R.T.
22.927	4.17 ng/ul	627322	Perylene-d12	24.204

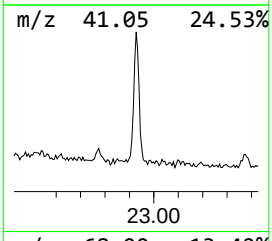
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Squalene	410	C30H50	000111-02-4	94
2	(2E,6E,10E)-3,7,11,15-Tetramethy...	318	C21H34O2	125456-63-5	90
3	Supraene	410	C30H50	007683-64-9	86
4	Squalene	410	C30H50	000111-02-4	83
5	3,7-Dimethyl-1-phenylsulfonyl-2,...	278	C16H22O2S	1000432-27-5	59

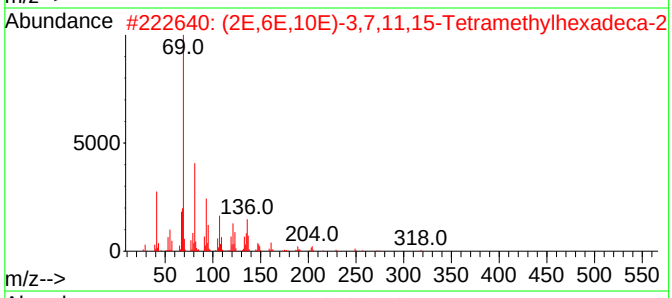


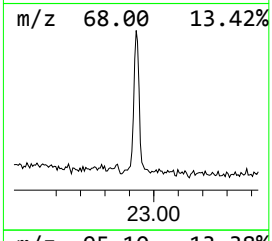


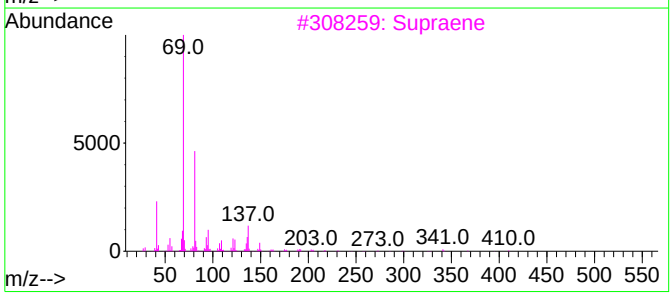


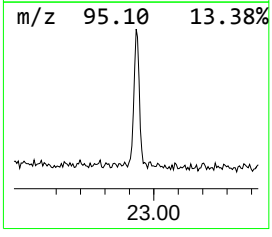












Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051023\
Data File : BP015043.D
Acq On : 10 May 2023 15:28
Operator : CG/JU
Sample : 02694-18
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
JREP8

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050323.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-Butanol, 3-me...	3.740	3.3	ng/ul	160827	1	8.146	983036	20.0
1-Propen-2-ol, ...	3.811	3.0	ng/ul	146240	1	8.146	983036	20.0
Butanoic acid	4.123	6.5	ng/ul	319172	1	8.146	983036	20.0
unknown-01	4.475	2.1	ng/ul	105596	1	8.146	983036	20.0
Crotonic acid	4.934	124.8	ng/ul	6136160	1	8.146	983036	20.0
2-Propanol, 1-p...	5.193	6.8	ng/ul	333058	1	8.146	983036	20.0
trans-2-Penteno...	6.234	5.4	ng/ul	266698	1	8.146	983036	20.0
2(3H)-Furanone,...	6.999	4.3	ng/ul	208730	1	8.146	983036	20.0
2-Propanol, 1-[...	7.728	3.2	ng/ul	158110	1	8.146	983036	20.0
2-Propanol, 1-(...	7.922	7.0	ng/ul	346024	1	8.146	983036	20.0
Phenylethyl Alc...	9.710	3.3	ng/ul	225602	2	10.981	1376500	20.0
1-Propanol, 2-(...	10.146	12.3	ng/ul	843343	2	10.981	1376500	20.0
4,5-Octanediol,...	11.752	3.6	ng/ul	247451	2	10.981	1376500	20.0
unknown-02	12.216	13.3	ng/ul	915483	2	10.981	1376500	20.0
unknown-03	12.681	285.4	ng/ul	19643500	2	10.981	1376500	20.0
2-Butenoic acid...	13.575	7.5	ng/ul	694314	3	14.787	1851020	20.0
3-Octanol, 2,2-...	13.740	7.8	ng/ul	723340	3	14.787	1851020	20.0
Ethanol, 2-[2-(...	14.322	5.7	ng/ul	523731	3	14.787	1851020	20.0
7-Methoxymethyl...	15.287	4.0	ng/ul	374889	3	14.787	1851020	20.0
Benzophenone	16.140	11.9	ng/ul	1101050	3	14.787	1851020	20.0
unknown-04	16.834	4.5	ng/ul	590975	4	17.540	2634060	20.0
unknown-05	17.045	115.0	ng/ul	15150700	4	17.540	2634060	20.0
unknown-06	17.740	2.6	ng/ul	345736	4	17.540	2634060	20.0
(E)-Hexadec-9-e...	18.287	2.8	ng/ul	373146	4	17.540	2634060	20.0
n-Hexadecanoic ...	18.404	16.9	ng/ul	2231490	4	17.540	2634060	20.0
unknown-07	18.528	2.6	ng/ul	347176	4	17.540	2634060	20.0
Octadecanoic acid	19.622	9.4	ng/ul	1302130	5	21.622	2759120	20.0
Behenic alcohol	21.292	6.4	ng/ul	877517	5	21.622	2759120	20.0
Squalene	22.927	4.2	ng/ul	627322	6	24.204	3006170	20.0