LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM020122\
Data File : BM033985.D
Acq On : 02 Feb 2022 21:59
Operator : CG/JU
Sample : N1355-09
Misc :
ALS Vial : 56 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0VF3

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_M\Methods\SFAM-EPA-BM013122.M
Title : SVOA CALIBRATION

Signal : TIC: BM033985.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.308	5	12	27	rVB	678999	858022	6.37%	1.100%
2	3.613	54	64	65	rBV3	45791	118922	0.88%	0.152%
3	3.649	65	70	80	rVB	837843	1130017	8.40%	1.449%
4	3.725	80	83	97	rVB	109043	178355	1.33%	0.229%
5	3.872	104	108	116	rBV	123869	166025	1.23%	0.213%
6	4.060	116	140	142	rBV3	279295	1324603	9.84%	1.698%
7	4.472	204	210	219	rVB	1154894	1579172	11.73%	2.025%
8	4.802	260	266	276	rBV	45389	79978	0.59%	0.103%
9	5.013	293	302	308	rBV	2032945	2927880	21.75%	3.754%
10	5.366	350	362	364	rVV4	28516	63951	0.48%	0.082%
11	5.407	364	369	376	rVV2	165469	262058	1.95%	0.336%
12	5.543	386	392	402	rVB	459762	860322	6.39%	1.103%
13	5.749	422	427	438	rVB	278283	385866	2.87%	0.495%
14	5.919	450	456	462	rVV	348647	548190	4.07%	0.703%
15	5.990	462	468	478	rVV	721069	1031312	7.66%	1.322%
16	6.390	529	536	540	rBV2	28847	67715	0.50%	0.087%
17	6.560	560	565	575	rVB	124651	211179	1.57%	0.271%
18	6.960	612	633	639	rBV4	140148	482211	3.58%	0.618%
19	7.007	639	641	646	rVV	62587	83558	0.62%	0.107%
20	7.084	646	654	661	rVV	599546	1066376	7.92%	1.367%
21	7.143	661	664	668	rVB	46822	62341	0.46%	0.080%
22	7.248	675	682	688	rVV	427918	670429	4.98%	0.860%
23	7.325	688	695	702	rVV3	68182	145406	1.08%	0.186%
24	7.466	707	719	724	rVV2	1114689	2426405	18.03%	3.111%
25	7.525	724	729	747	rVB	1061297	1910181	14.19%	2.449%
26	7.719	755	762	775	rVB	251710	513010	3.81%	0.658%
27	7.925	789	797	803	rBV	524173	867497	6.45%	1.112%
28	7.995	803	809	813	rBV	63615	101062	0.75%	0.130%
29	8.043	813	817	826	rVB	57429	98977	0.74%	0.127%
30	8.190	826	842	856	rBV	6871184	13459249	100.00%	17.257%
31	8.466	878	889	896	rBV5	35520	101664	0.76%	0.130%
32	8.631	909	917	925	rVV	719203	1262088	9.38%	1.618%
33	8.748	932	937	945	rVB	82672	146374	1.09%	0.188%
34	8.878	954	959	964	rBV2	43088	72836	0.54%	0.093%
35	8.937	964	969	978	rVB3	33995	63218	0.47%	0.081%
36	9.037	978	986	989	rBV3	82001	156978	1.17%	0.201%

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Misc :
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Instrument : BNA_M
ClientSampleId : C0VF3

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_M\Methods\SFAM-EPA-BM013122.M
Title : SVOA CALIBRATION

37	9.084	989	994	1000	rVV	543041	860661	6.39%	1.104%
38	9.166	1000	1008	1013	rVV	206225	288284	2.14%	0.370%
39	9.242	1013	1021	1027	rBV	168991	334702	2.49%	0.429%
40	9.619	1082	1085	1094	rVV2	35101	74366	0.55%	0.095%
41	9.813	1109	1118	1124	rBV	476376	837628	6.22%	1.074%
42	10.095	1139	1166	1172	rBV	575424	2721880	20.22%	3.490%
43	10.354	1202	1210	1219	rVV2	768436	1500847	11.15%	1.924%
44	10.513	1231	1237	1245	rVV	233936	366013	2.72%	0.469%
45	10.731	1265	1274	1280	rVV	702528	1328708	9.87%	1.704%
46	10.789	1280	1284	1291	rVB3	51856	96960	0.72%	0.124%
47	10.866	1291	1297	1310	rBV	302002	570596	4.24%	0.732%
48	11.089	1330	1335	1343	rBV2	43247	83508	0.62%	0.107%
49	11.507	1400	1406	1417	rBV2	65454	164381	1.22%	0.211%
50	11.642	1424	1429	1437	rBV2	52126	101175	0.75%	0.130%
51	12.201	1521	1524	1529	rVB	46171	70358	0.52%	0.090%
52	12.260	1529	1534	1539	rVB	54016	85449	0.63%	0.110%
53	12.501	1568	1575	1589	rBV2	34258	116489	0.87%	0.149%
54	12.978	1650	1656	1666	rBV7	35897	100208	0.74%	0.128%
55	13.389	1719	1726	1733	rBV3	120171	233733	1.74%	0.300%
56	13.607	1758	1763	1771	rVB	62368	98139	0.73%	0.126%
57	13.972	1819	1825	1830	rVB	1084083	1524337	11.33%	1.954%
58	14.066	1837	1841	1845	rBV	68497	97276	0.72%	0.125%
59	14.107	1845	1848	1855	rVB	100632	144660	1.07%	0.185%
60	14.254	1867	1873	1879	rVV	1281774	1803682	13.40%	2.313%
61	14.466	1904	1909	1918	rBV	82921	107239	0.80%	0.137%
62	14.560	1918	1925	1931	rBV	976704	1450833	10.78%	1.860%
63	14.748	1952	1957	1969	rBV	356850	567533	4.22%	0.728%
64	15.077	2007	2013	2022	rVB4	31179	68822	0.51%	0.088%
65	15.413	2065	2070	2075	rVB	75790	95205	0.71%	0.122%
66	15.548	2084	2093	2100	rBV	1567925	2165845	16.09%	2.777%
67	15.660	2106	2112	2120	rBV	479889	651473	4.84%	0.835%
68	16.095	2180	2186	2189	rBV2	39651	62614	0.47%	0.080%
69	16.236	2205	2210	2213	rBV	101624	146635	1.09%	0.188%
70	16.442	2241	2245	2251	rBV2	38481	60685	0.45%	0.078%
71	16.618	2266	2275	2278	rBV5	45900	102767	0.76%	0.132%
72	17.060	2346	2350	2356	rVV	76239	120165	0.89%	0.154%
73	17.295	2384	2390	2396	rBV	1063450	1445120	10.74%	1.853%
74	17.395	2400	2407	2418	rVB	1479008	2070504	15.38%	2.655%
75	17.801	2472	2476	2479	rBV	48981	62239	0.46%	0.080%
76	18.189	2538	2542	2558	rBV2	203731	383407	2.85%	0.492%

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Misc :
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Instrument :
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C0VF3

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % of largest Peak
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Stop Thrs : 0 Peak Location: TOP

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Title : SVOA CALIBRATION

77	18.489	2589	2593	2604	rBV2	57530	102761	0.76%	0.132%
78	18.918	2659	2666	2676	rVB10	38215	112855	0.84%	0.145%
79	19.124	2696	2701	2706	rBV4	45371	99354	0.74%	0.127%
80	19.189	2706	2712	2717	rVB6	69678	126644	0.94%	0.162%
81	19.295	2724	2730	2733	rVV3	132686	244127	1.81%	0.313%
82	19.348	2736	2739	2746	rVV	125153	207468	1.54%	0.266%
83	19.465	2756	2759	2762	rBV	56415	62878	0.47%	0.081%
84	19.612	2780	2784	2788	rVB2	58801	74181	0.55%	0.095%
85	19.677	2790	2795	2800	rVV	1604455	2229709	16.57%	2.859%
86	19.718	2800	2802	2807	rVV	127254	169263	1.26%	0.217%
87	19.771	2808	2811	2815	rVV5	59175	88498	0.66%	0.113%
88	19.818	2815	2819	2825	rVB	142309	178433	1.33%	0.229%
89	20.542	2936	2942	2945	rBV3	95801	171027	1.27%	0.219%
90	20.942	3006	3010	3014	rBV6	58266	110130	0.82%	0.141%
91	21.171	3044	3049	3059	rBV2	296914	443447	3.29%	0.569%
92	21.459	3093	3098	3103	rBV	1654158	2130707	15.83%	2.732%
93	22.083	3201	3204	3209	rVB4	131537	184765	1.37%	0.237%
94	22.153	3213	3216	3222	rVB7	145132	197808	1.47%	0.254%
95	23.306	3408	3412	3416	rBV	190306	286649	2.13%	0.368%
96	23.371	3416	3423	3433	rVB	4260572	7253288	53.89%	9.300%
97	23.694	3471	3478	3484	rBV2	1047388	2038725	15.15%	2.614%
98	23.847	3498	3504	3514	rVB2	738527	1484418	11.03%	1.903%
99	24.624	3633	3636	3644	rVB	229000	455574	3.38%	0.584%
100	24.847	3669	3674	3685	rVB	404667	891296	6.62%	1.143%

Sum of corrected areas: 77992558

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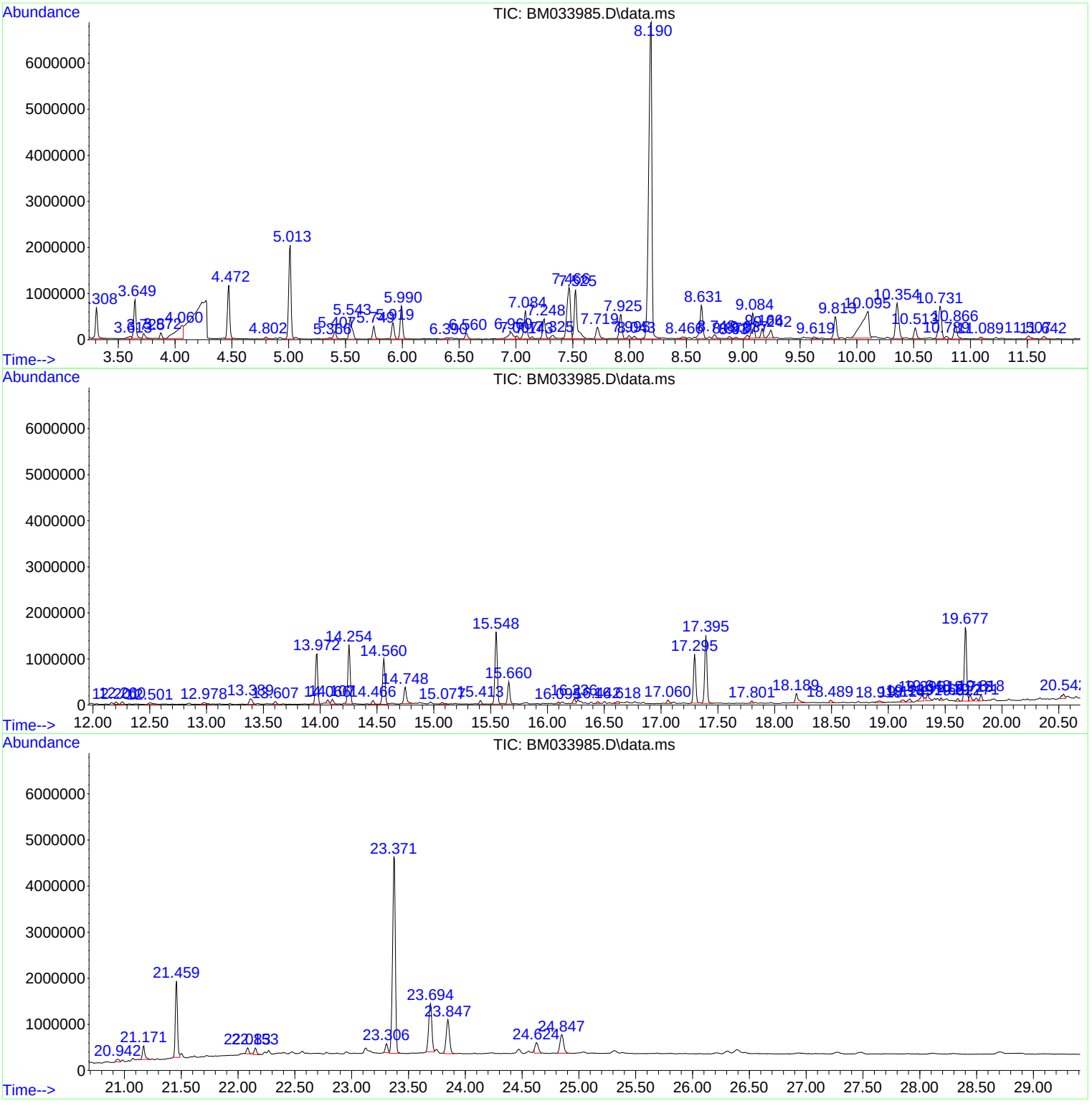
ALS Vial : 56 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0VF3

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM013122.M
Quant Title : SVOA CALIBRATION

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



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ALS Vial : 56 Sample Multiplier: 1

Instrument :
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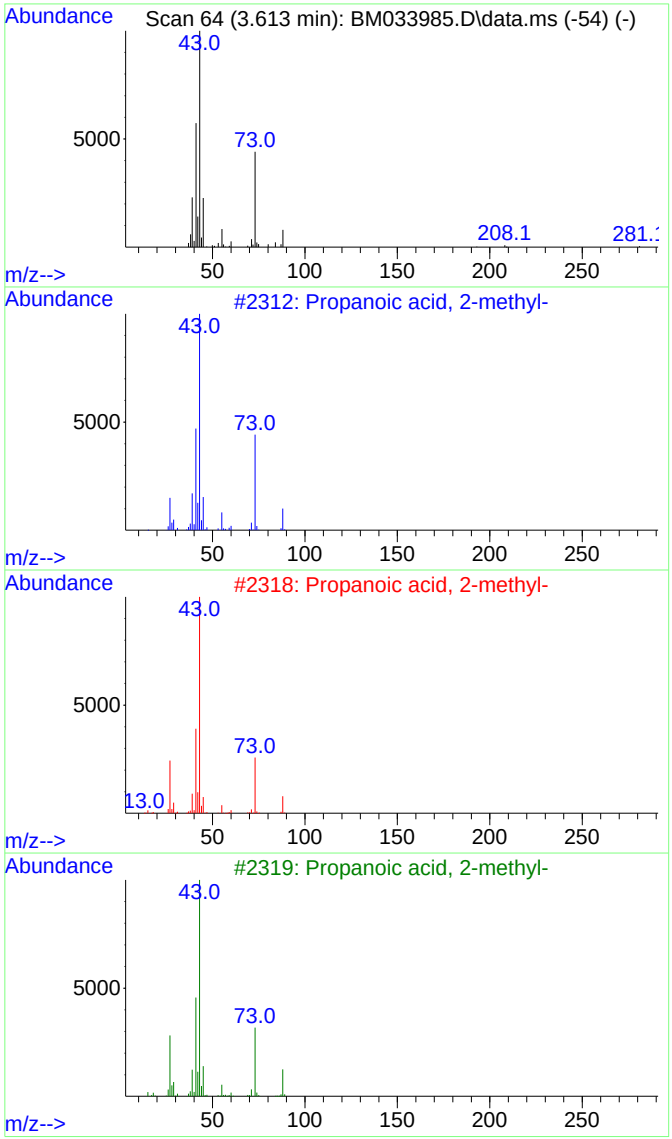
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM013122.M
Quant Title : SVOA CALIBRATION

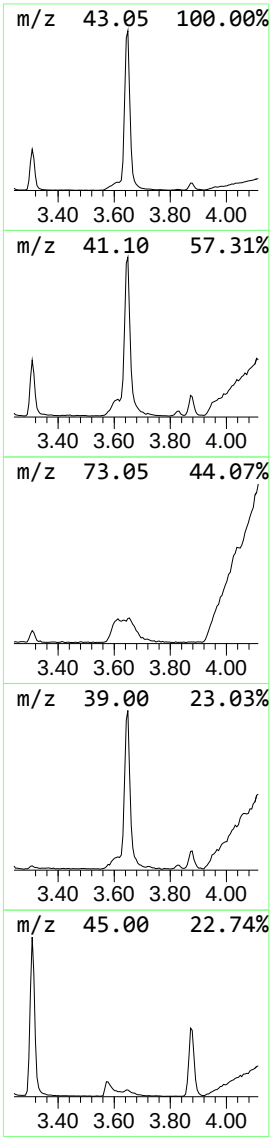
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 Propanoic acid, 2-methyl- Concentration Rank 30

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.613	2.74 ng/ul	118922	1,4-Dichlorobenzene-d4	7.925

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Propanoic acid, 2-methyl-	88	C4H8O2	000079-31-2	72
2	Propanoic acid, 2-methyl-	88	C4H8O2	000079-31-2	59
3	Propanoic acid, 2-methyl-	88	C4H8O2	000079-31-2	53
4	Propanoic acid, 2-methyl-	88	C4H8O2	000079-31-2	53
5	Propanoic acid, 2-methyl-	88	C4H8O2	000079-31-2	50





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ALS Vial : 56 Sample Multiplier: 1

Instrument :
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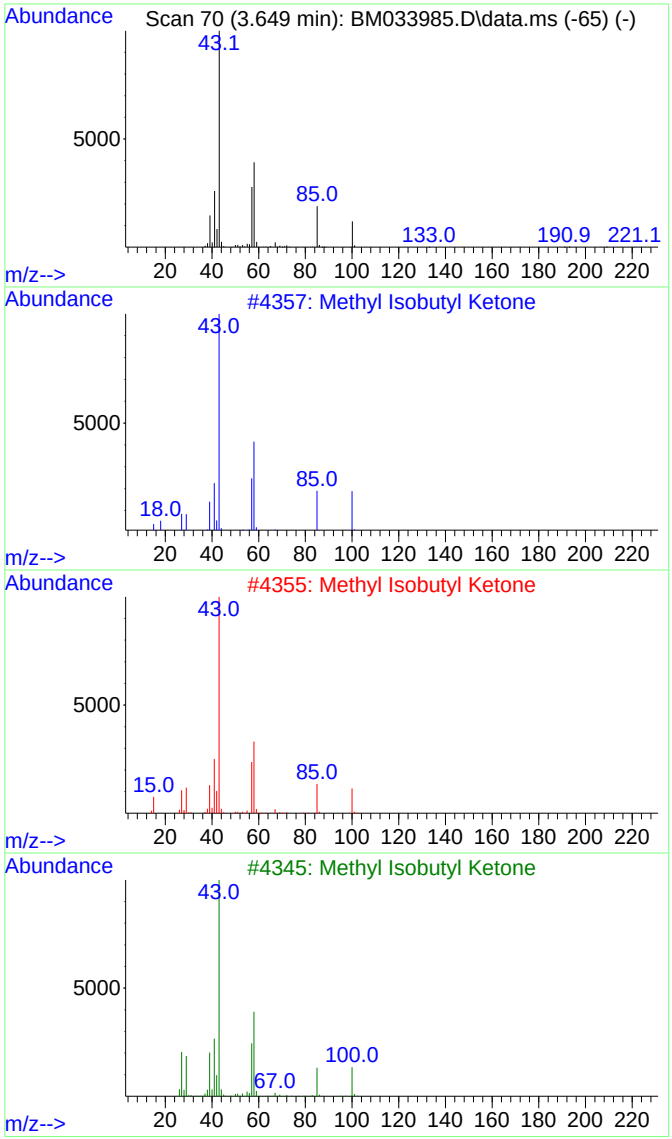
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM013122.M
Quant Title : SVOA CALIBRATION

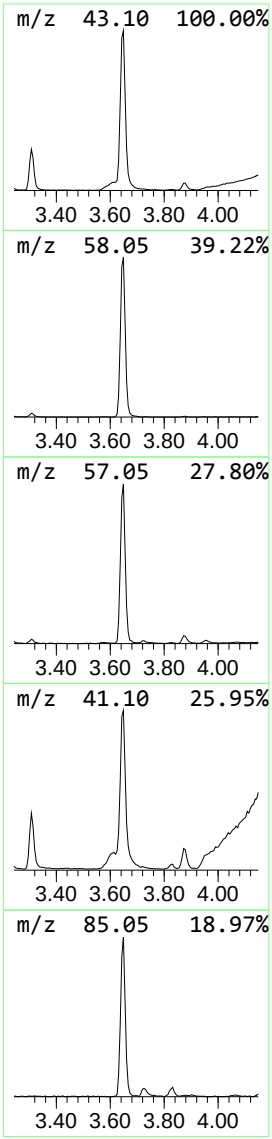
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 Methyl Isobutyl Ketone Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.649	26.05 ng/ul	1130020	1,4-Dichlorobenzene-d4	7.925

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Methyl Isobutyl Ketone	100	C6H12O	000108-10-1	91
2	Methyl Isobutyl Ketone	100	C6H12O	000108-10-1	87
3	Methyl Isobutyl Ketone	100	C6H12O	000108-10-1	86
4	Methyl Isobutyl Ketone	100	C6H12O	000108-10-1	86
5	2-Hexanone	100	C6H12O	000591-78-6	78





Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM020122\
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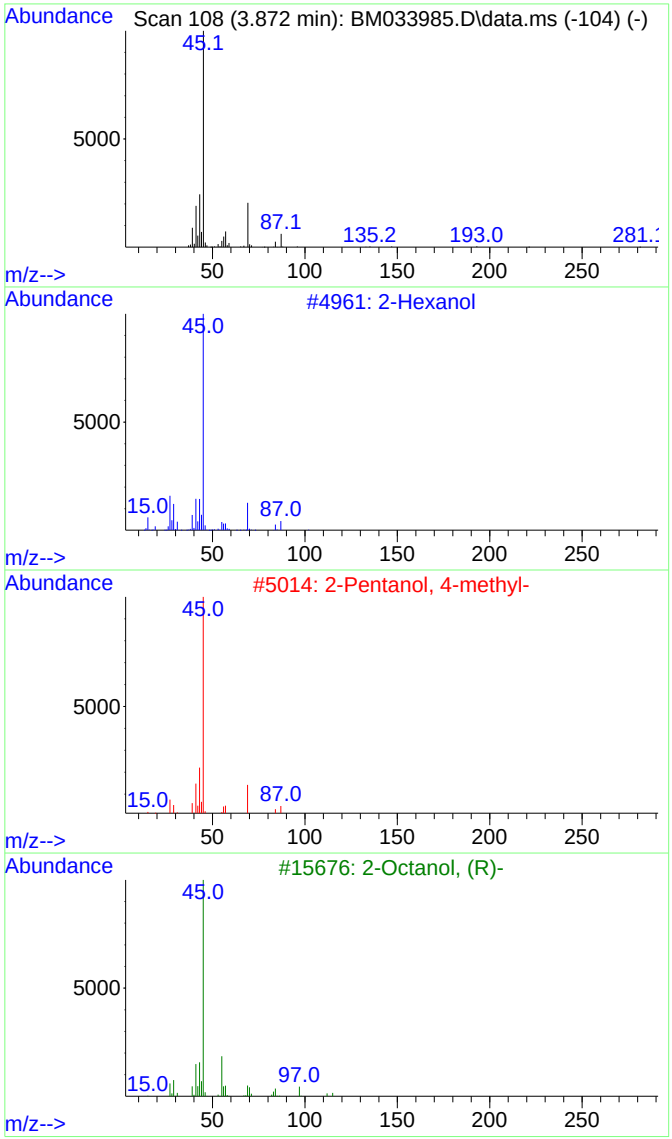
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM013122.M
Quant Title : SVOA CALIBRATION

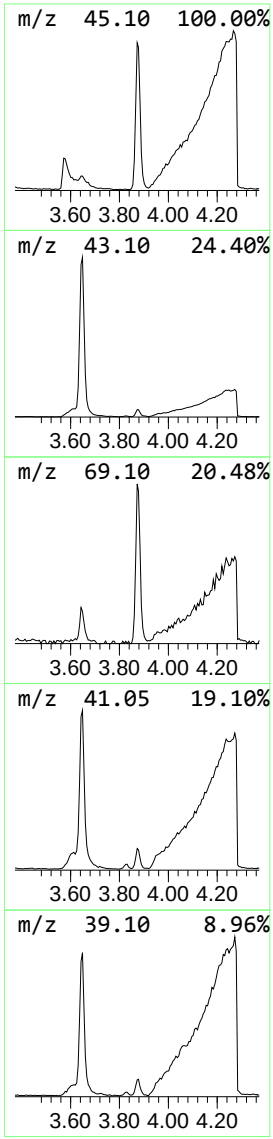
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 2-Hexanol Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.872	3.83 ng/ul	166025	1,4-Dichlorobenzene-d4	7.925

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Hexanol	102	C6H14O	000626-93-7	83
2	2-Pentanol, 4-methyl-	102	C6H14O	000108-11-2	78
3	2-Octanol, (R)-	130	C8H18O	005978-70-1	78
4	2-Octanol	130	C8H18O	000123-96-6	78
5	2-Hexanol	102	C6H14O	000626-93-7	74





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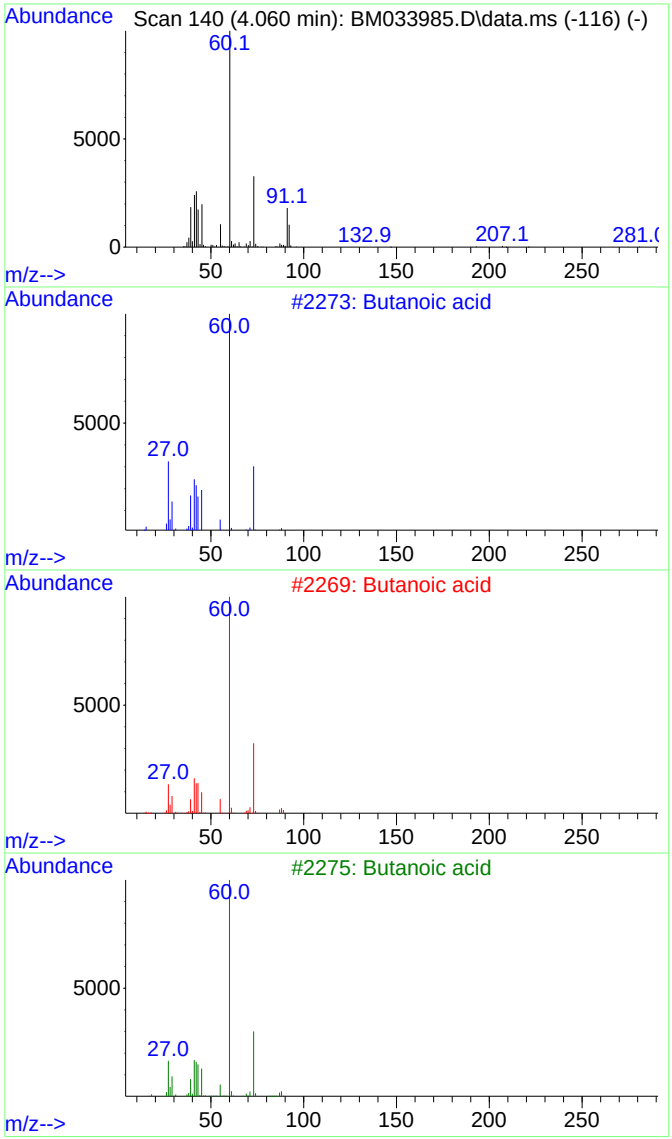
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM013122.M
Quant Title : SVOA CALIBRATION

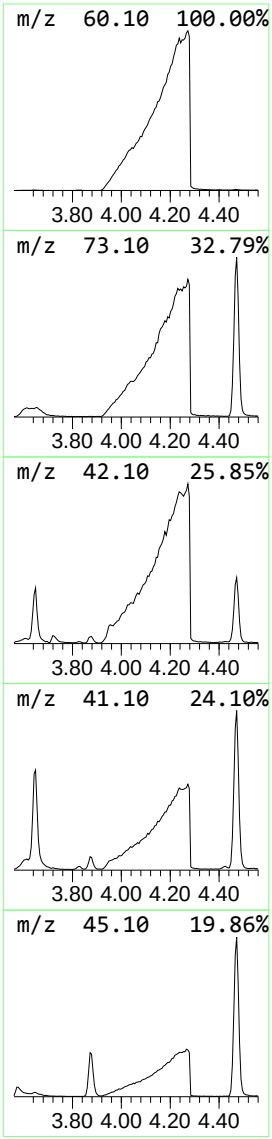
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Butanoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.060	30.54 ng/ul	1324600	1,4-Dichlorobenzene-d4	7.925

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butanoic acid	88	C4H8O2	000107-92-6	64
2	Butanoic acid	88	C4H8O2	000107-92-6	58
3	Butanoic acid	88	C4H8O2	000107-92-6	53
4	Butanoic acid	88	C4H8O2	000107-92-6	53
5	Silver butanoate	194	C4H7AgO2	005076-24-4	38





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Misc :
ALS Vial : 56 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0VF3

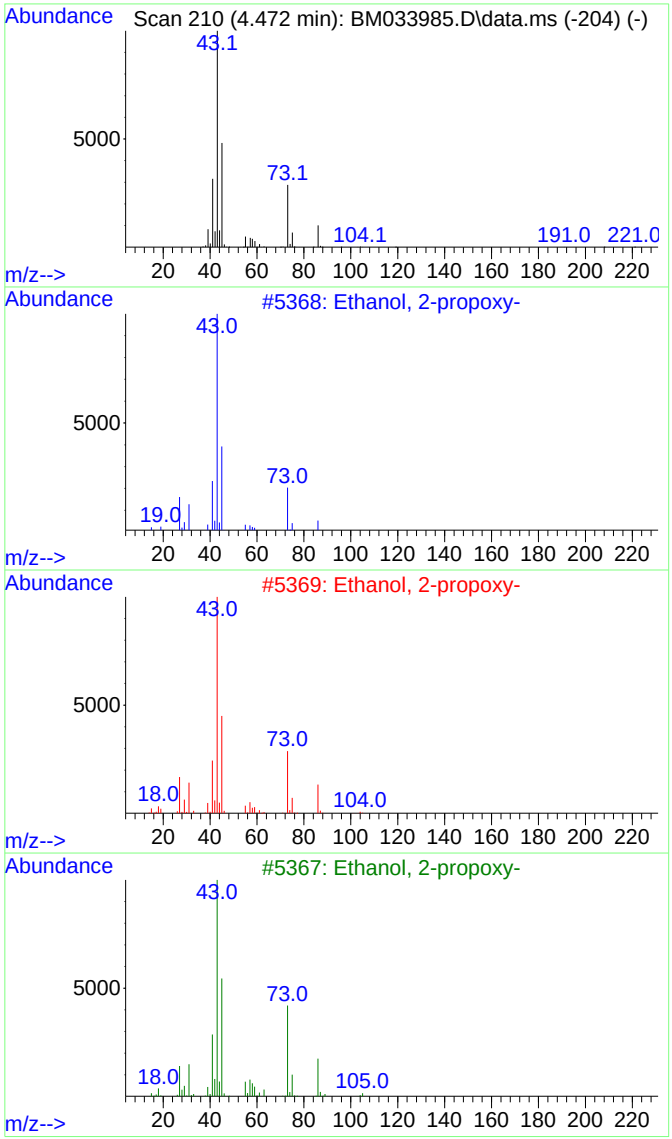
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM013122.M
Quant Title : SVOA CALIBRATION

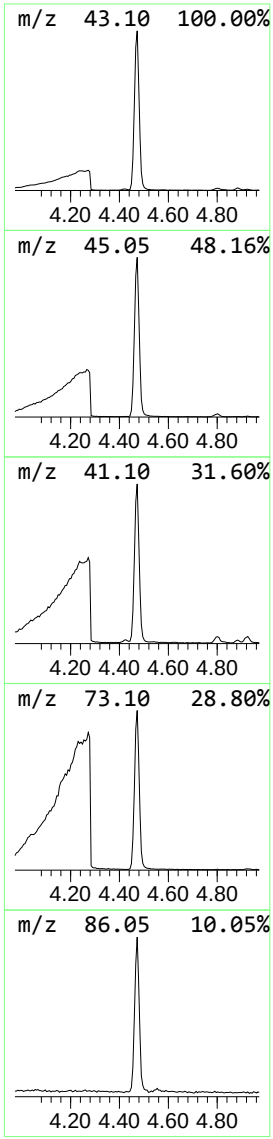
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Ethanol, 2-propoxy- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.472	36.41 ng/ul	1579170	1,4-Dichlorobenzene-d4	7.925

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethanol, 2-propoxy-	104	C5H12O2	002807-30-9	83
2	Ethanol, 2-propoxy-	104	C5H12O2	002807-30-9	50
3	Ethanol, 2-propoxy-	104	C5H12O2	002807-30-9	33
4	Ethanol, 2-methoxy-, acetate	118	C5H10O3	000110-49-6	33
5	Acetic acid ethenyl ester	86	C4H6O2	000108-05-4	9





Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM020122\
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Acq On : 02 Feb 2022 21:59
Operator : CG/JU
Sample : N1355-09
Misc :
ALS Vial : 56 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0VF3

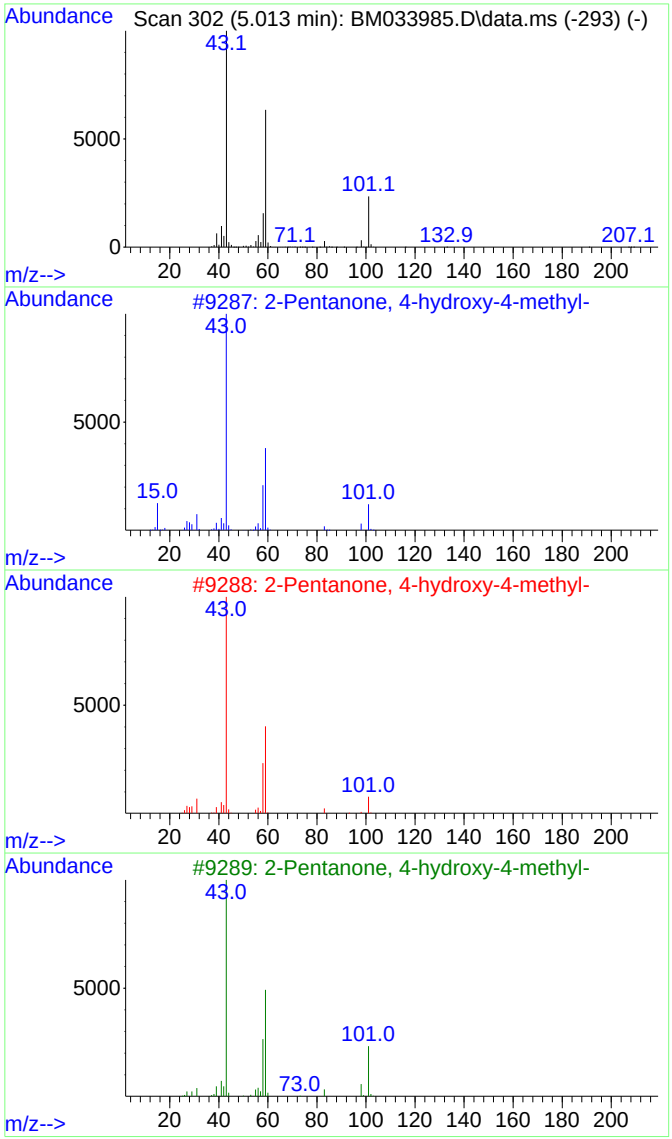
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM013122.M
Quant Title : SVOA CALIBRATION

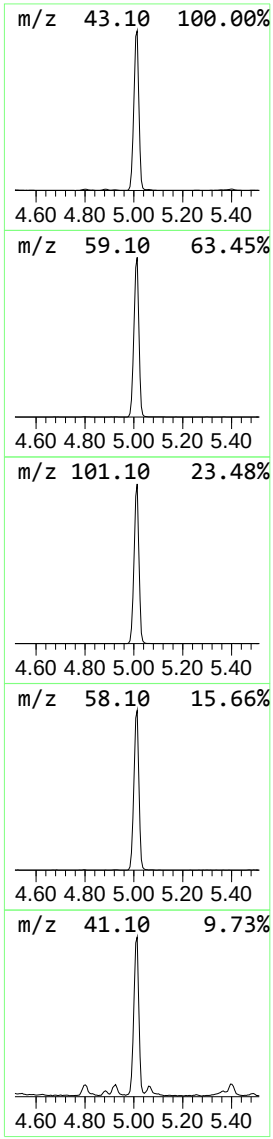
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.013	67.50 ng/ul	2927880	1,4-Dichlorobenzene-d4	7.925

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
3	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	40
4	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	28
5	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	23





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Sample : N1355-09
Misc :
ALS Vial : 56 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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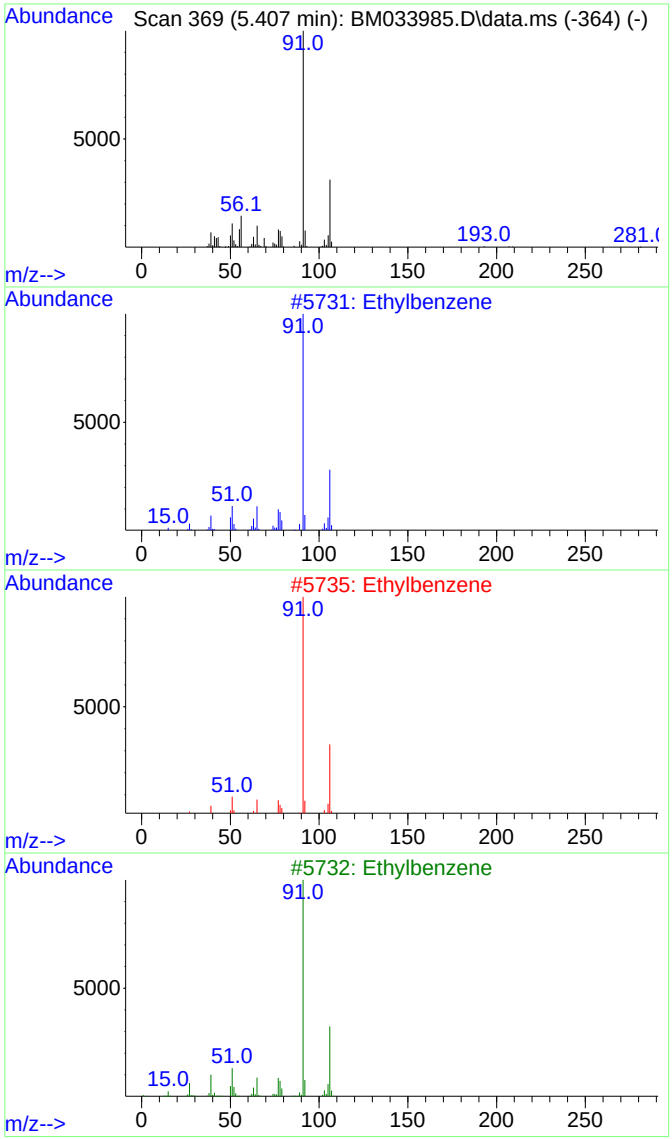
Quant Title : SVOA CALIBRATION

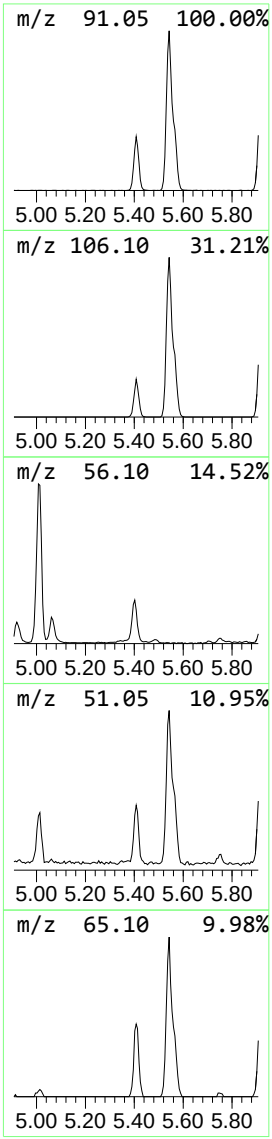
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 Ethylbenzene Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.407	6.04 ng/ul	262058	1,4-Dichlorobenzene-d4	7.925

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethylbenzene	106	C8H10	000100-41-4	94
2	Ethylbenzene	106	C8H10	000100-41-4	93
3	Ethylbenzene	106	C8H10	000100-41-4	93
4	Ethylbenzene	106	C8H10	000100-41-4	93
5	o-Xylene	106	C8H10	000095-47-6	72





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ALS Vial : 56 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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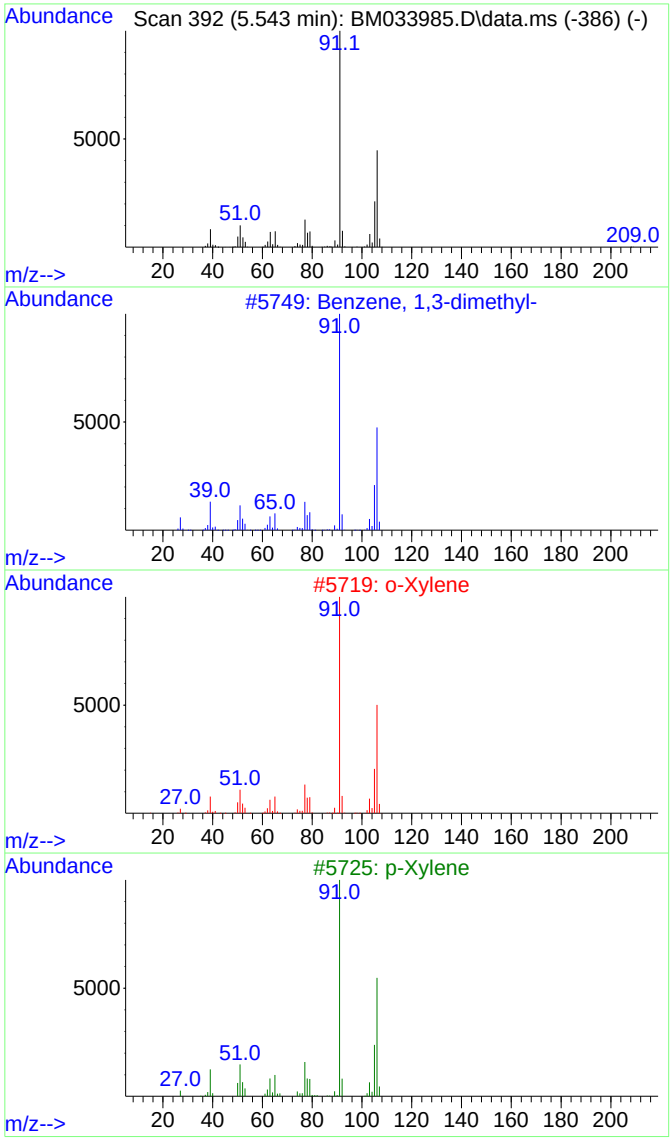
Quant Title : SVOA CALIBRATION

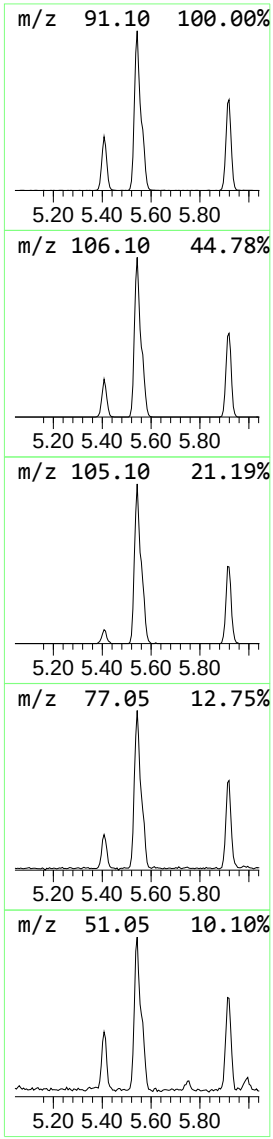
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 Benzene, 1,3-dimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.543	19.83 ng/ul	860322	1,4-Dichlorobenzene-d4	7.925

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,3-dimethyl-	106	C8H10	000108-38-3	97
2	o-Xylene	106	C8H10	000095-47-6	97
3	p-Xylene	106	C8H10	000106-42-3	97
4	p-Xylene	106	C8H10	000106-42-3	97
5	Benzene, 1,3-dimethyl-	106	C8H10	000108-38-3	97





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ClientSampleId :
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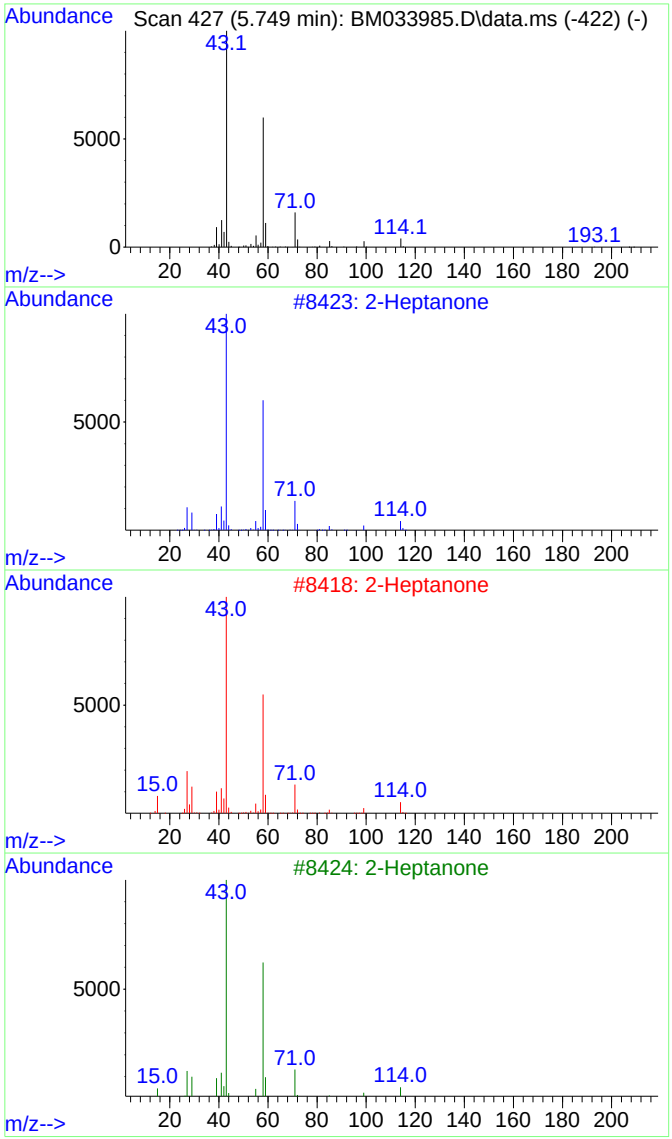
Quant Title : SVOA CALIBRATION

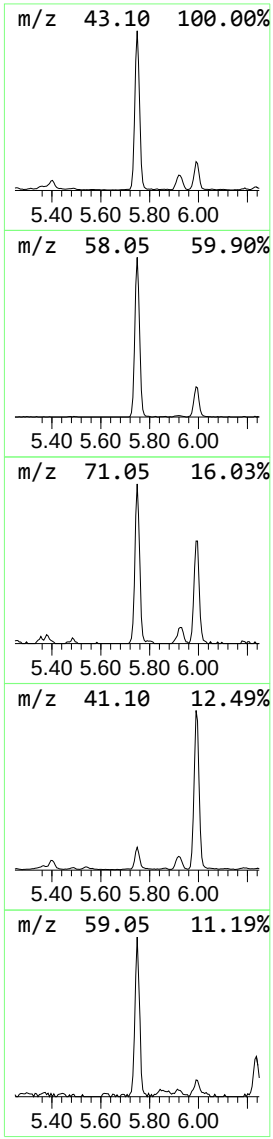
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 2-Heptanone Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.749	8.90 ng/ul	385866	1,4-Dichlorobenzene-d4	7.925

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Heptanone	114	C7H14O	000110-43-0	91
2	2-Heptanone	114	C7H14O	000110-43-0	91
3	2-Heptanone	114	C7H14O	000110-43-0	90
4	2-Hexanone, 5-methyl-	114	C7H14O	000110-12-3	86
5	2-Heptanone	114	C7H14O	000110-43-0	83





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Misc :
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Instrument :
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

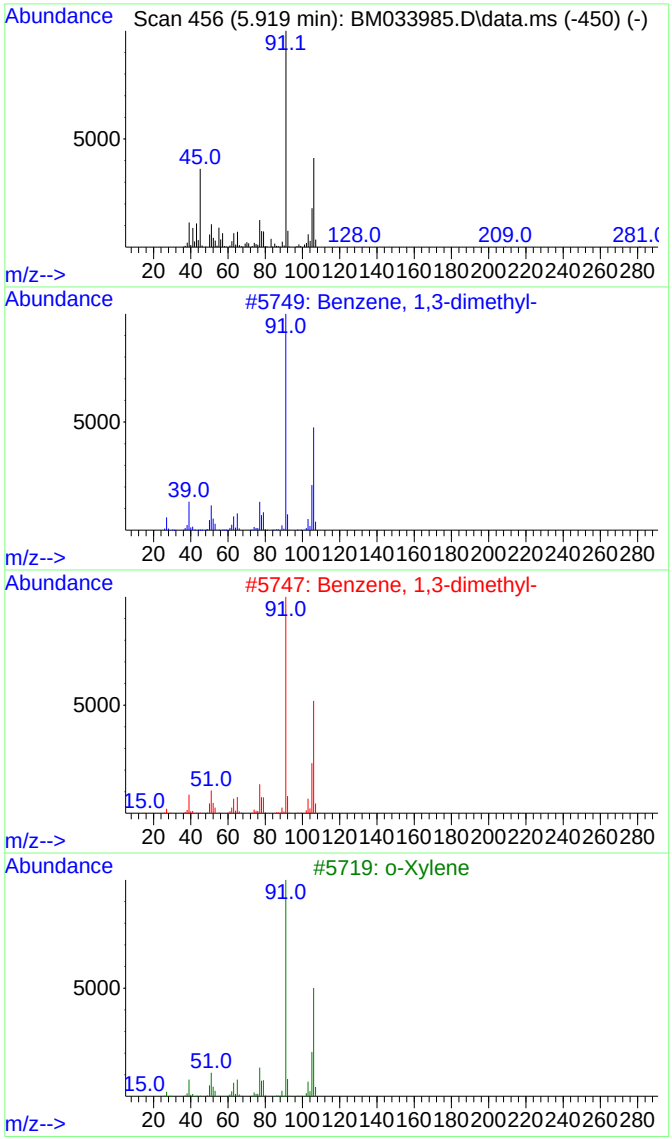
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

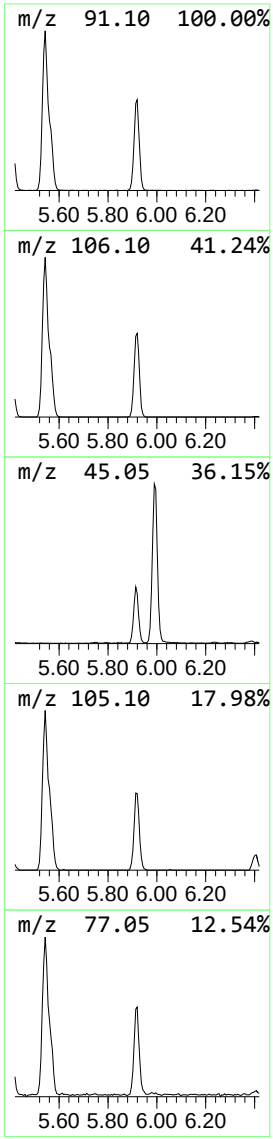
****

Peak Number 10 o-Xylene Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.919	12.64 ng/ul	548190	1,4-Dichlorobenzene-d4	7.925

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,3-dimethyl-	106	C8H10	000108-38-3	95
2	Benzene, 1,3-dimethyl-	106	C8H10	000108-38-3	95
3	o-Xylene	106	C8H10	000095-47-6	95
4	o-Xylene	106	C8H10	000095-47-6	95
5	p-Xylene	106	C8H10	000106-42-3	95





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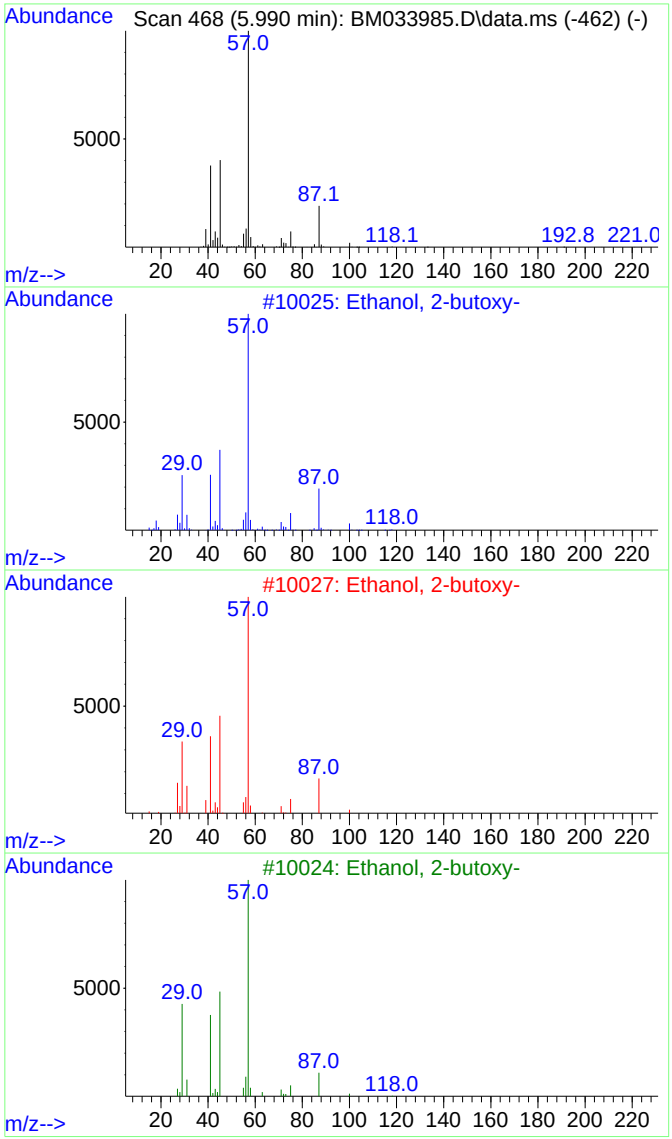
Quant Title : SVOA CALIBRATION

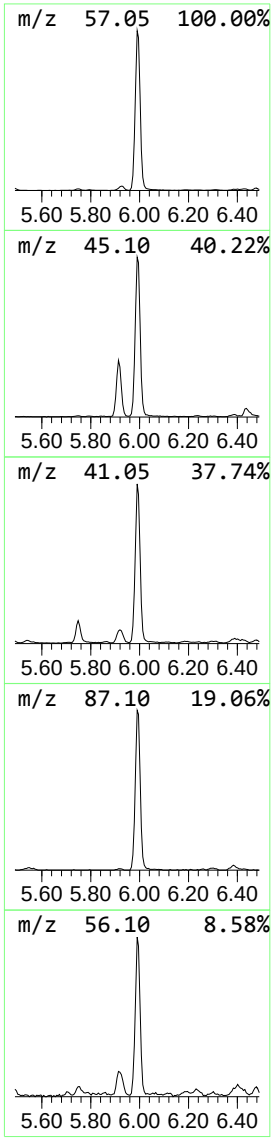
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 11 Ethanol, 2-butoxy- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.990	23.78 ng/ul	1031310	1,4-Dichlorobenzene-d4	7.925

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethanol, 2-butoxy-	118	C6H14O2	000111-76-2	91
2	Ethanol, 2-butoxy-	118	C6H14O2	000111-76-2	83
3	Ethanol, 2-butoxy-	118	C6H14O2	000111-76-2	64
4	Ethanol, 2-butoxy-	118	C6H14O2	000111-76-2	49
5	Ethylene glycol monoisobutyl ether	118	C6H14O2	004439-24-1	47





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Misc :
ALS Vial : 56 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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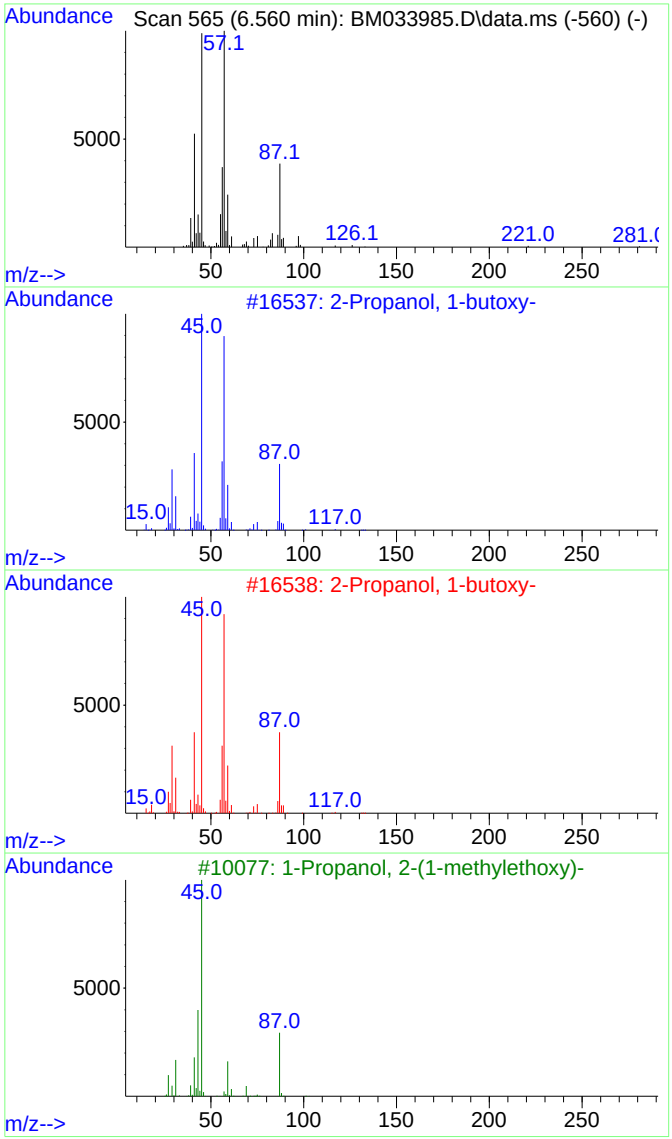
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM013122.M
Quant Title : SVOA CALIBRATION

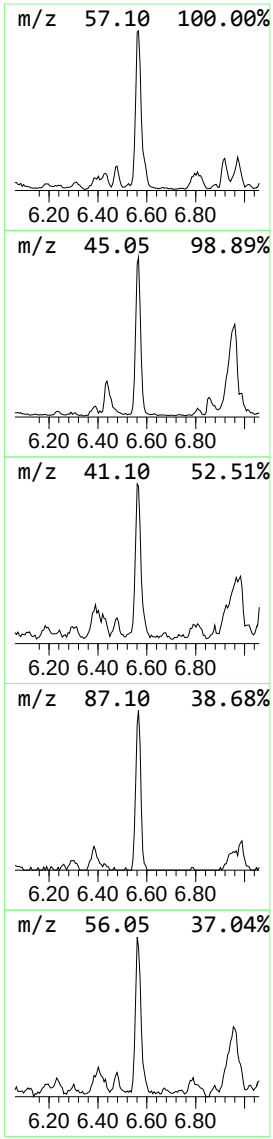
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 12 2-Propanol, 1-butoxy- Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.560	4.87 ng/ul	211179	1,4-Dichlorobenzene-d4	7.925

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Propanol, 1-butoxy-	132	C7H16O2	005131-66-8	90
2	2-Propanol, 1-butoxy-	132	C7H16O2	005131-66-8	90
3	1-Propanol, 2-(1-methylethoxy)-	118	C6H14O2	003944-37-4	53
4	Butane, 1-(1-methylpropoxy)-	130	C8H18O	000999-65-5	47
5	3,3-Dimethylbutane-2-ol	102	C6H14O	000464-07-3	42





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ClientSampleId :
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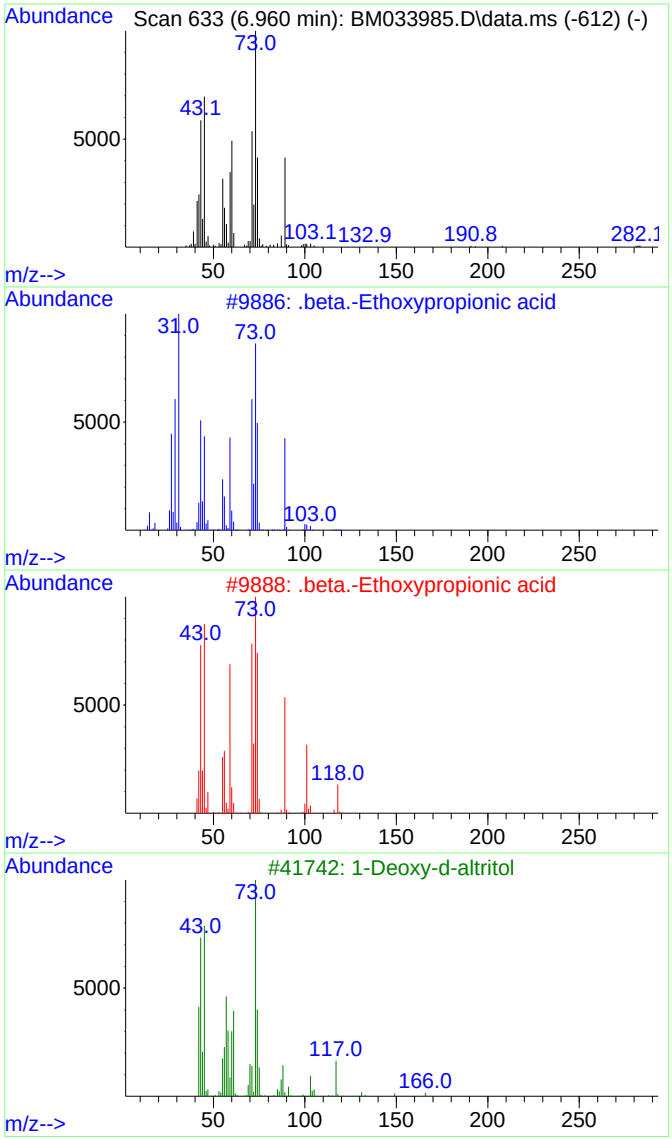
Quant Title : SVOA CALIBRATION

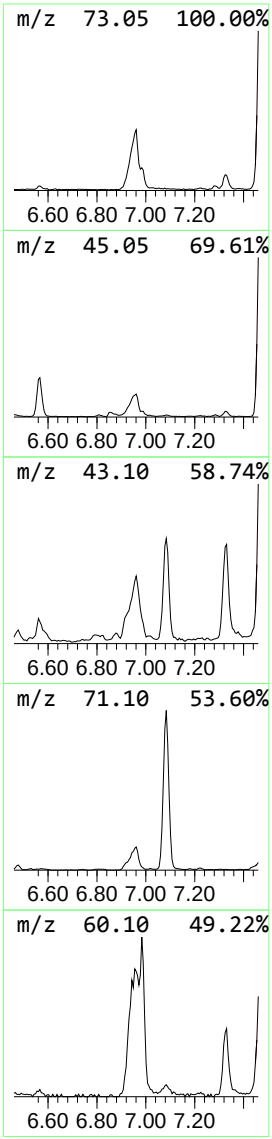
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 13 .beta.-Ethoxypropionic acid Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.960	11.12 ng/ul	482211	1,4-Dichlorobenzene-d4	7.925

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	.beta.-Ethoxypropionic acid	118	C5H10O3	004324-38-3	58
2	.beta.-Ethoxypropionic acid	118	C5H10O3	004324-38-3	53
3	1-Deoxy-d-altritol	166	C6H14O5	068832-18-8	32
4	Acetaldehyde ethyl isoamyl acetal	160	C9H20O2	1000430-98-8	27
5	1-Propanol, 3-methoxy-2-methyl-	104	C5H12O2	051866-93-4	25





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Instrument :
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ClientSampleId :
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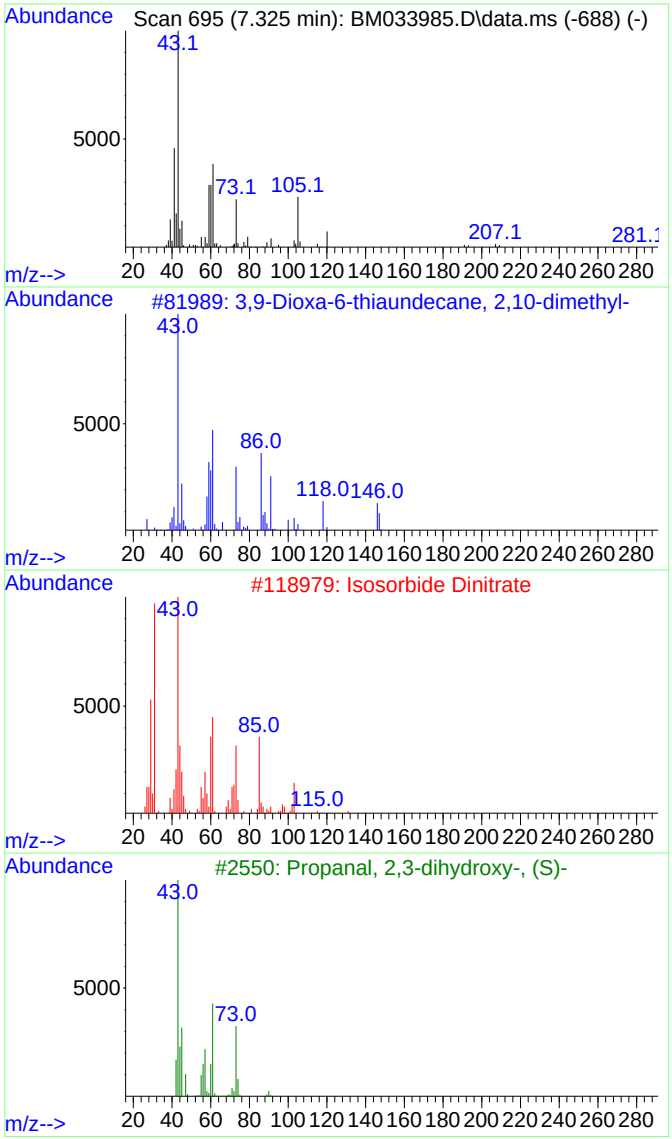
Quant Title : SVOA CALIBRATION

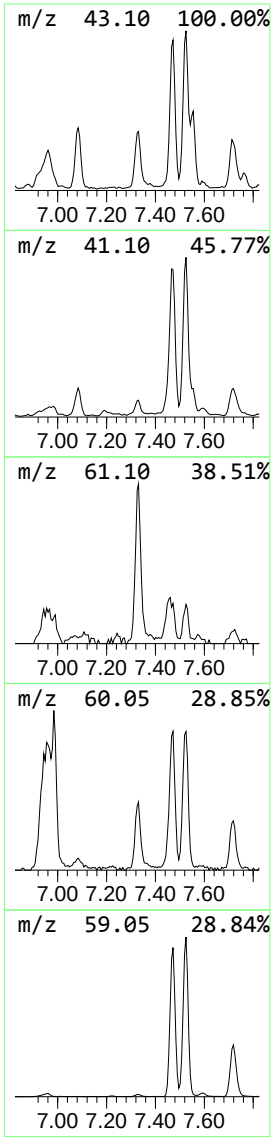
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 14 unknown-01 Concentration Rank 27

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.325	3.35 ng/ul	145406	1,4-Dichlorobenzene-d4	7.925

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3,9-Dioxa-6-thiaundecane, 2,10-d...	206	C10H22O2S	097916-00-2	36
2	Isosorbide Dinitrate	236	C6H8N2O8	000087-33-2	33
3	Propanal, 2,3-dihydroxy-, (S)-	90	C3H6O3	000497-09-6	28
4	n-Propyl acetate	102	C5H10O2	000109-60-4	25
5	Diacetamide	101	C4H7NO2	000625-77-4	25





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ClientSampleId :
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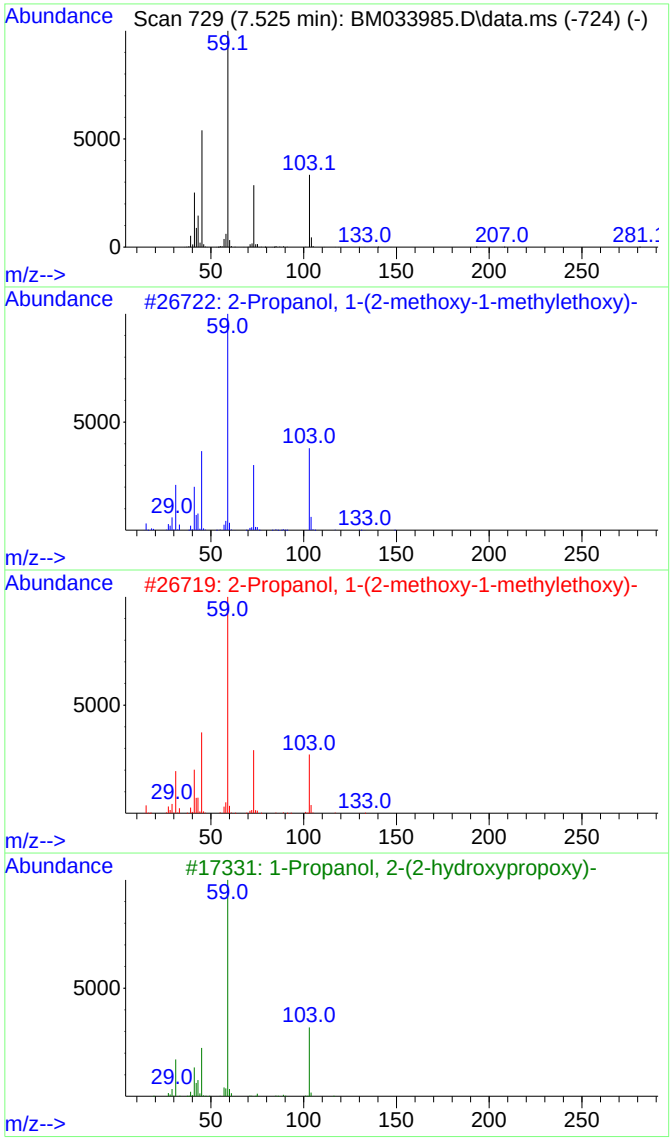
Quant Title : SVOA CALIBRATION

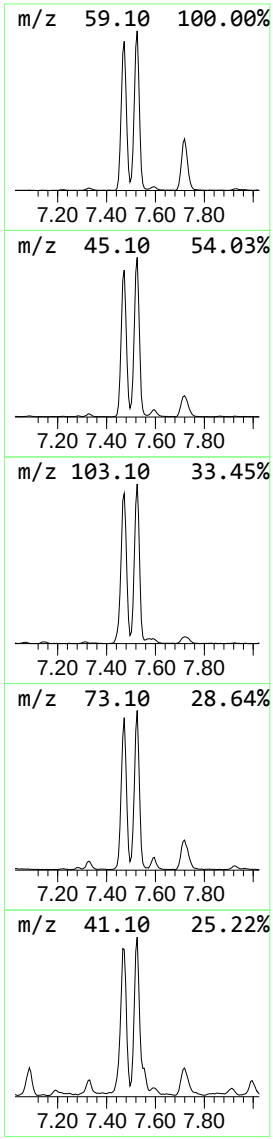
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 15 2-Propanol, 1-(2-methoxy-1-... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.525	44.04 ng/ul	1910180	1,4-Dichlorobenzene-d4	7.925

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Propanol, 1-(2-methoxy-1-methy...	148	C7H16O3	020324-32-7	80
2	2-Propanol, 1-(2-methoxy-1-methy...	148	C7H16O3	020324-32-7	72
3	1-Propanol, 2-(2-hydroxypropoxy)-	134	C6H14O3	000106-62-7	53
4	Dipropylene glycol (isomer 3)	134	C6H14O3	1000506-25-5	50
5	1-Propanol, 2,2'-oxybis-	134	C6H14O3	000108-61-2	50





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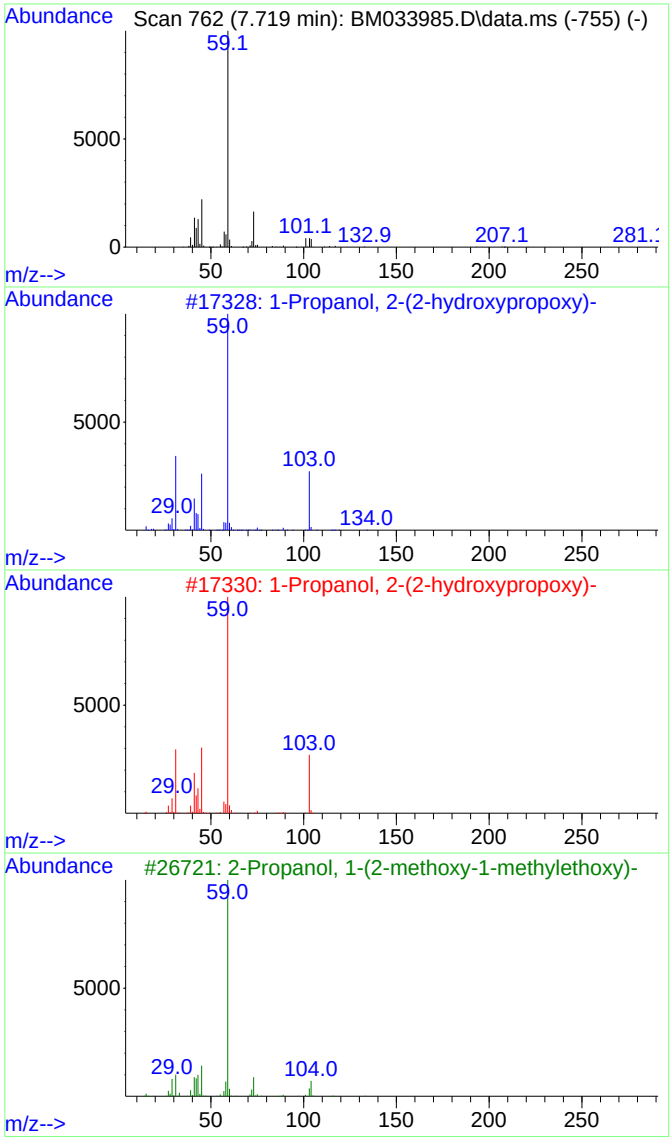
Quant Title : SVOA CALIBRATION

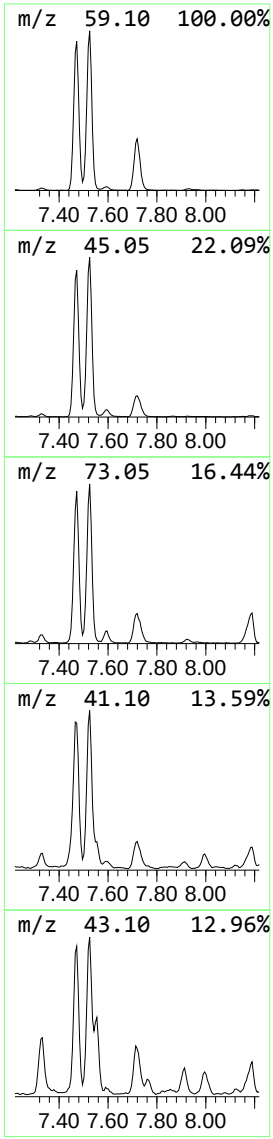
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 16 1-Propanol, 2-(2-hydroxypro... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.719	11.83 ng/ul	513010	1,4-Dichlorobenzene-d4	7.925

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Propanol, 2-(2-hydroxypropoxy)-	134	C6H14O3	000106-62-7	64
2	1-Propanol, 2-(2-hydroxypropoxy)-	134	C6H14O3	000106-62-7	64
3	2-Propanol, 1-(2-methoxy-1-methy...	148	C7H16O3	020324-32-7	64
4	1-Propanol, 2,2'-oxybis-	134	C6H14O3	000108-61-2	64
5	2-Propanol, 1-(2-methoxypropoxy)-	148	C7H16O3	013429-07-7	64





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Sample : N1355-09
Misc :
ALS Vial : 56 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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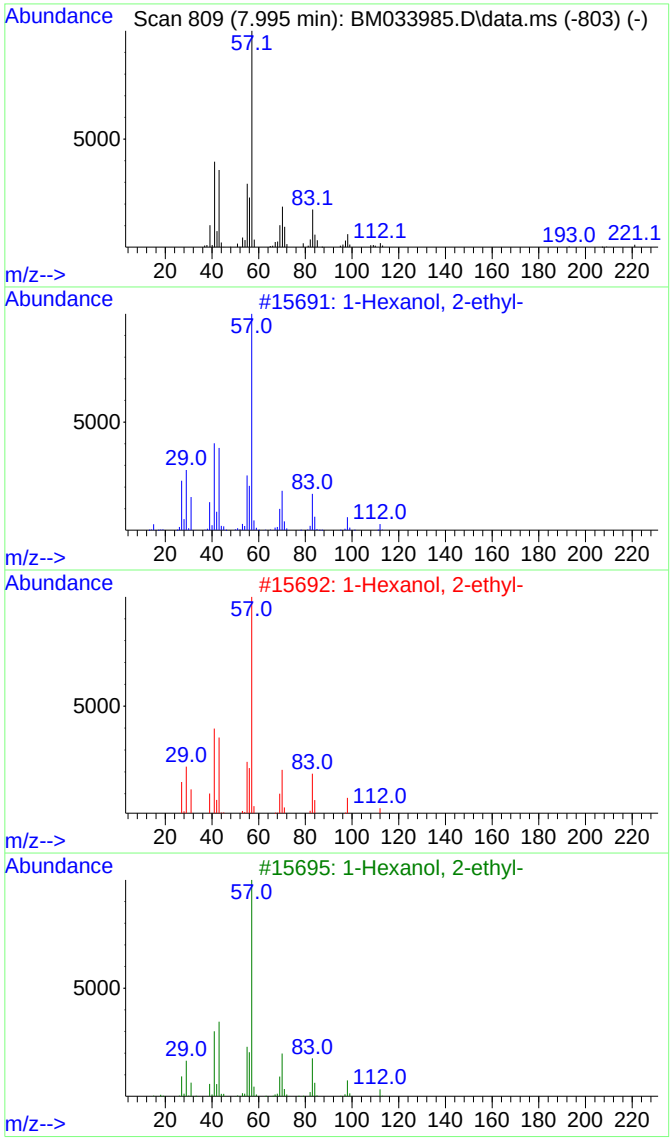
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM013122.M
Quant Title : SVOA CALIBRATION

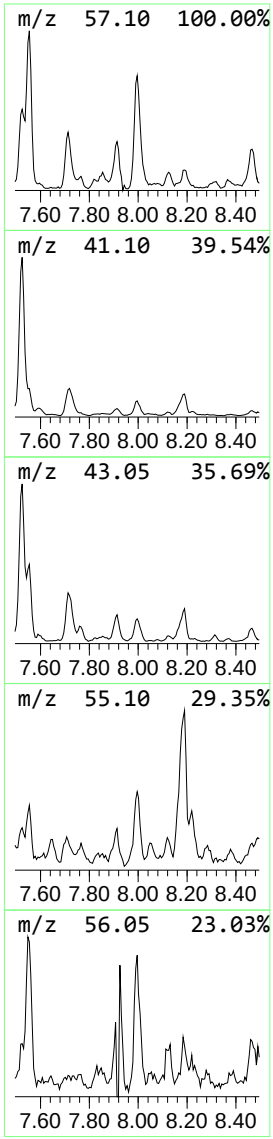
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 17 1-Hexanol, 2-ethyl- Concentration Rank 33

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.995	2.33 ng/ul	101062	1,4-Dichlorobenzene-d4	7.925

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	78
2	1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	72
3	1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	72
4	1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	59
5	1-Isobutoxy-2-ethylhexane	186	C12H26O	1000139-90-3	59





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Acq On : 02 Feb 2022 21:59
Operator : CG/JU
Sample : N1355-09
Misc :
ALS Vial : 56 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0VF3

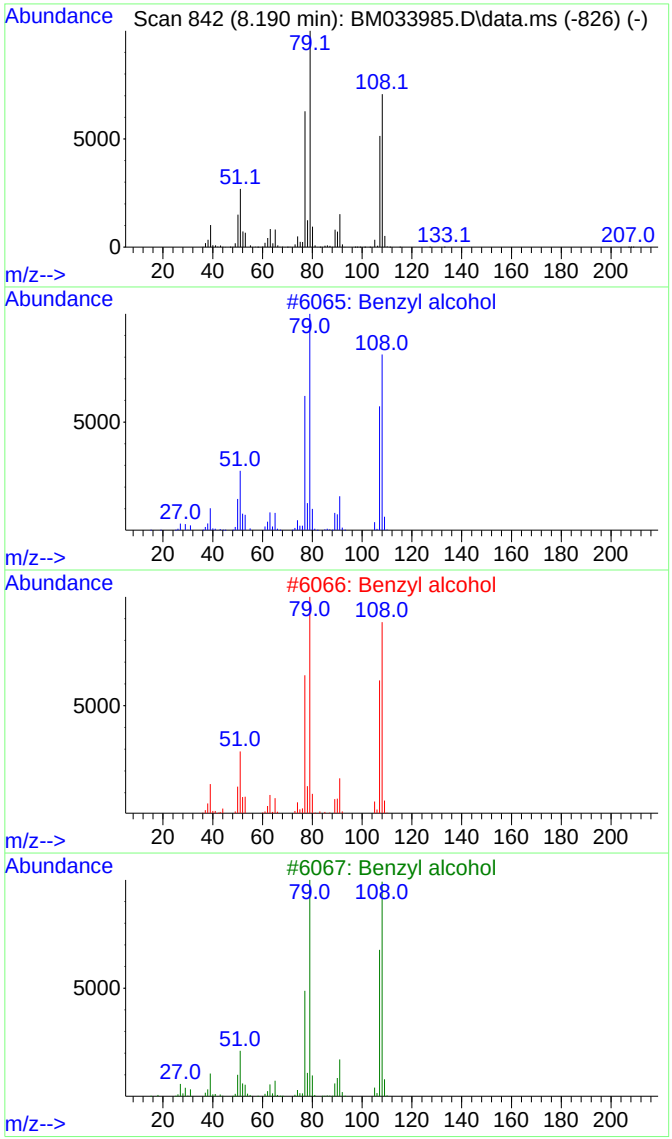
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM013122.M
Quant Title : SVOA CALIBRATION

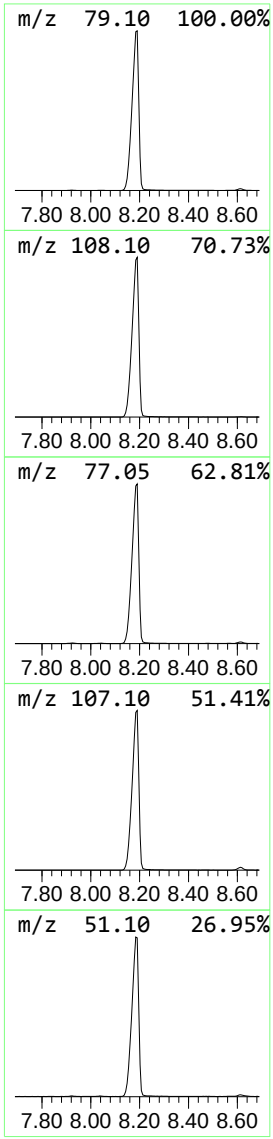
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 19 Benzyl alcohol Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.190	310.30 ng/ul	13459200	1,4-Dichlorobenzene-d4	7.925

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzyl alcohol	108	C7H8O	000100-51-6	98
2	Benzyl alcohol	108	C7H8O	000100-51-6	97
3	Benzyl alcohol	108	C7H8O	000100-51-6	97
4	Benzyl alcohol	108	C7H8O	000100-51-6	94
5	N-Cbz-glycylglycine p-nitropheny...	387	C18H17N3O7	013574-81-7	91





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Instrument :
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ClientSampleId :
C0VF3

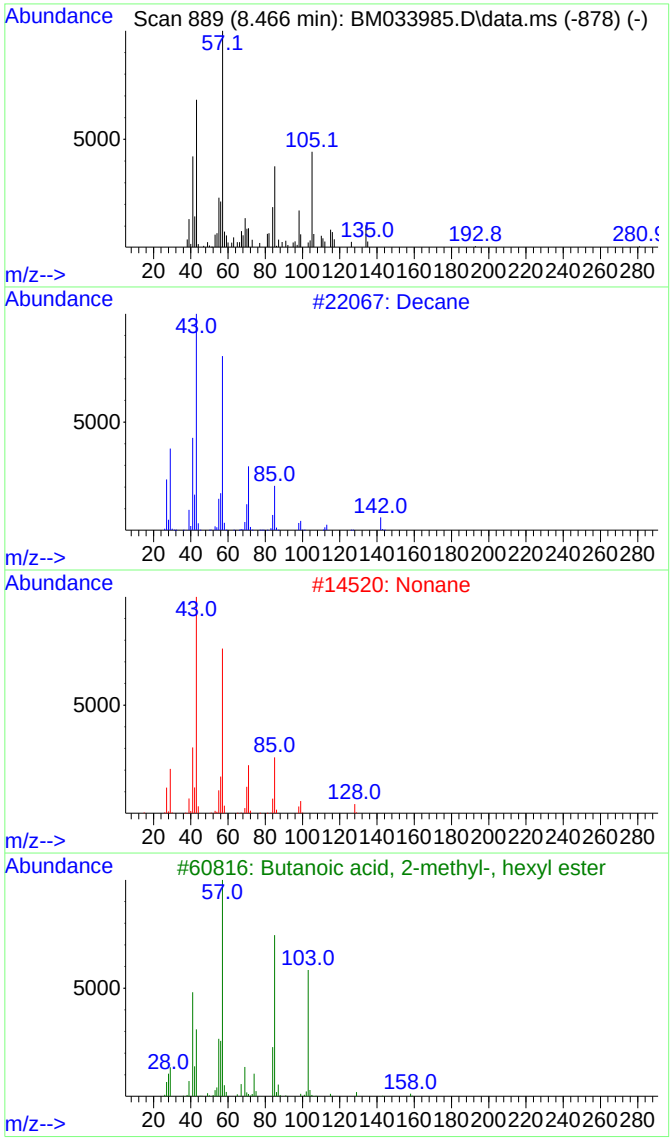
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM013122.M
Quant Title : SVOA CALIBRATION

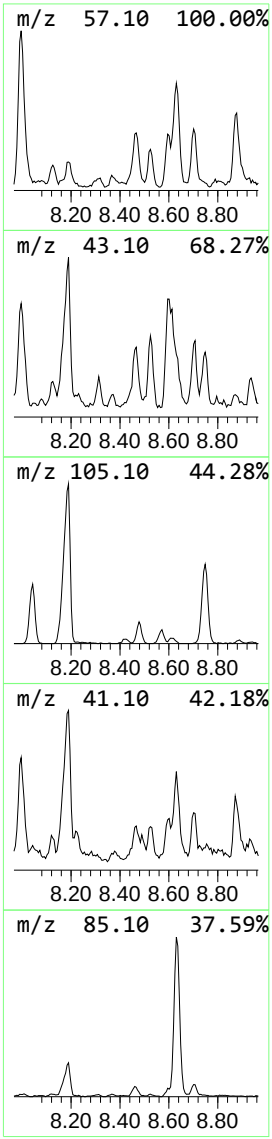
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 20 (DEL) Alkane: Straight-Chai... Concentration Rank 32

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.466	2.34 ng/ul	101664	1,4-Dichlorobenzene-d4	7.925

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Decane	142	C10H22	000124-18-5	50
2	Nonane	128	C9H20	000111-84-2	50
3	Butanoic acid, 2-methyl-, hexyl ...	186	C11H22O2	010032-15-2	47
4	Butanoic acid, 3-oxo-, 1,1-dimet...	158	C8H14O3	001694-31-1	47
5	Dodecane	170	C12H26	000112-40-3	47





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ALS Vial : 56 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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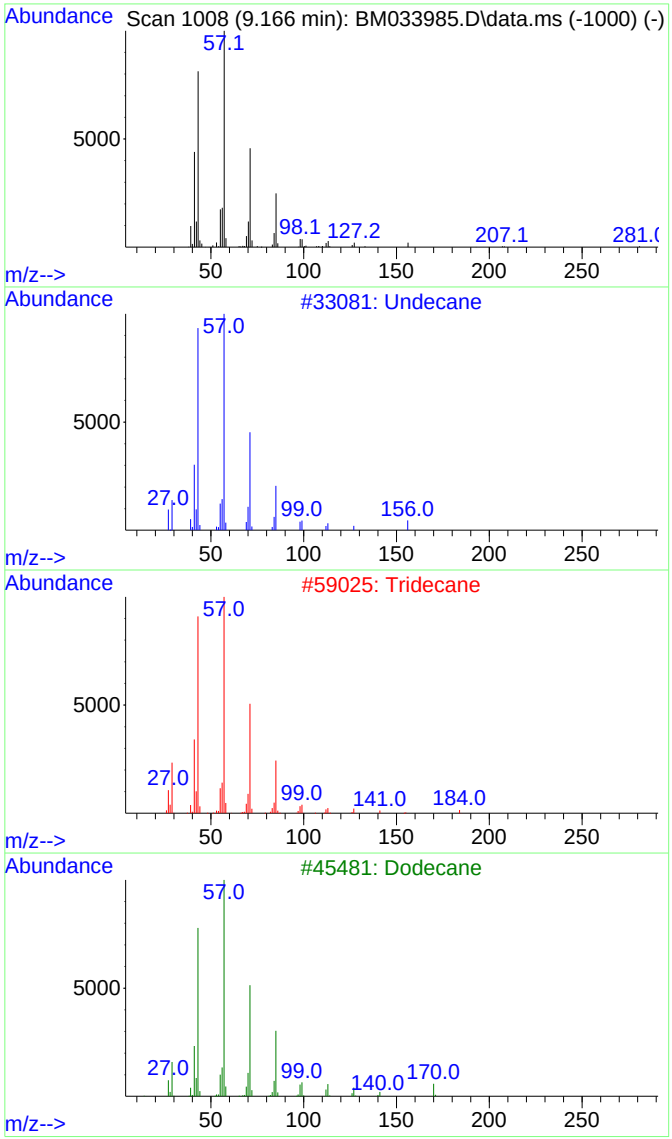
Quant Title : SVOA CALIBRATION

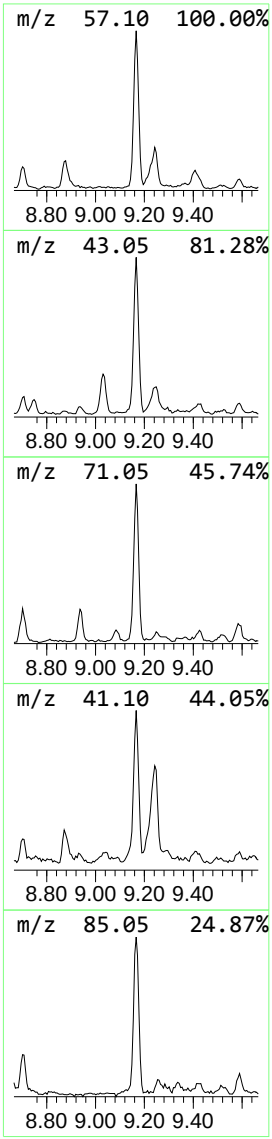
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 21 (DEL) Alkane: Straight-Chai... Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.166	6.65 ng/ul	288284	1,4-Dichlorobenzene-d4	7.925

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Undecane	156	C11H24	001120-21-4	91
2	Tridecane	184	C13H28	000629-50-5	86
3	Dodecane	170	C12H26	000112-40-3	83
4	Carbonic acid, prop-1-en-2-yl tr...	284	C17H32O3	1000382-90-8	83
5	Hexadecane	226	C16H34	000544-76-3	78





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ALS Vial : 56 Sample Multiplier: 1

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ClientSampleId :
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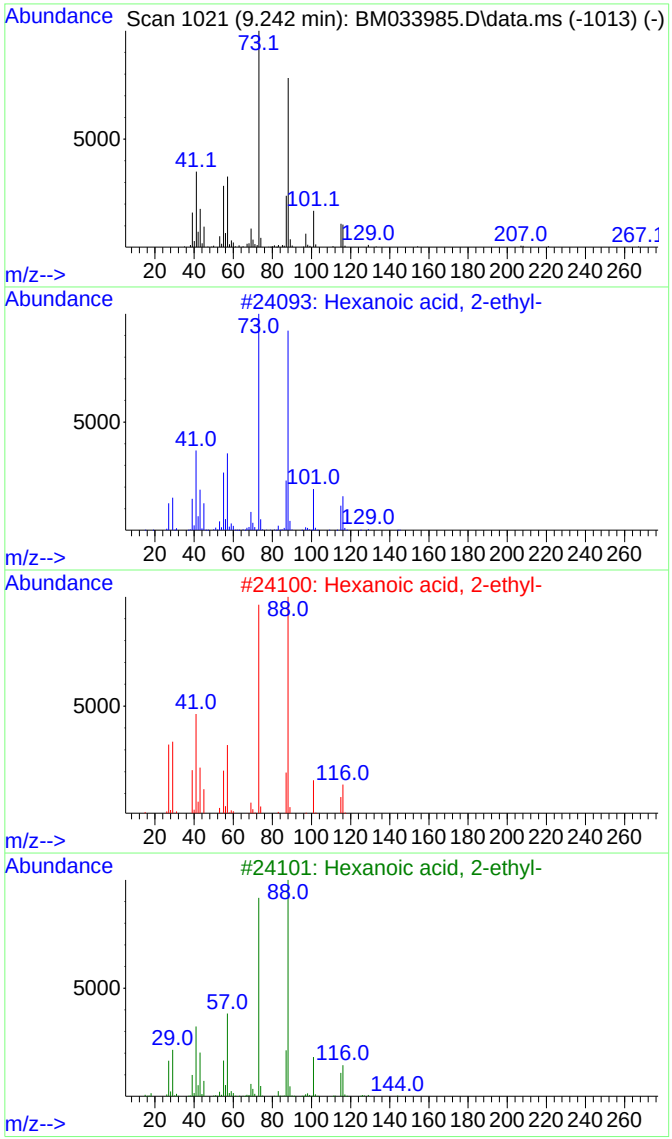
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM013122.M
Quant Title : SVOA CALIBRATION

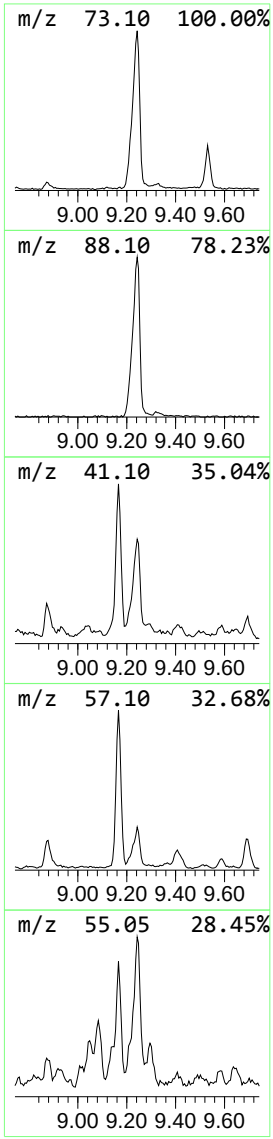
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 22 Hexanoic acid, 2-ethyl- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.242	7.72 ng/ul	334702	1,4-Dichlorobenzene-d4	7.925

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	86
2	Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	86
3	Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	80
4	Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	78
5	Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	78





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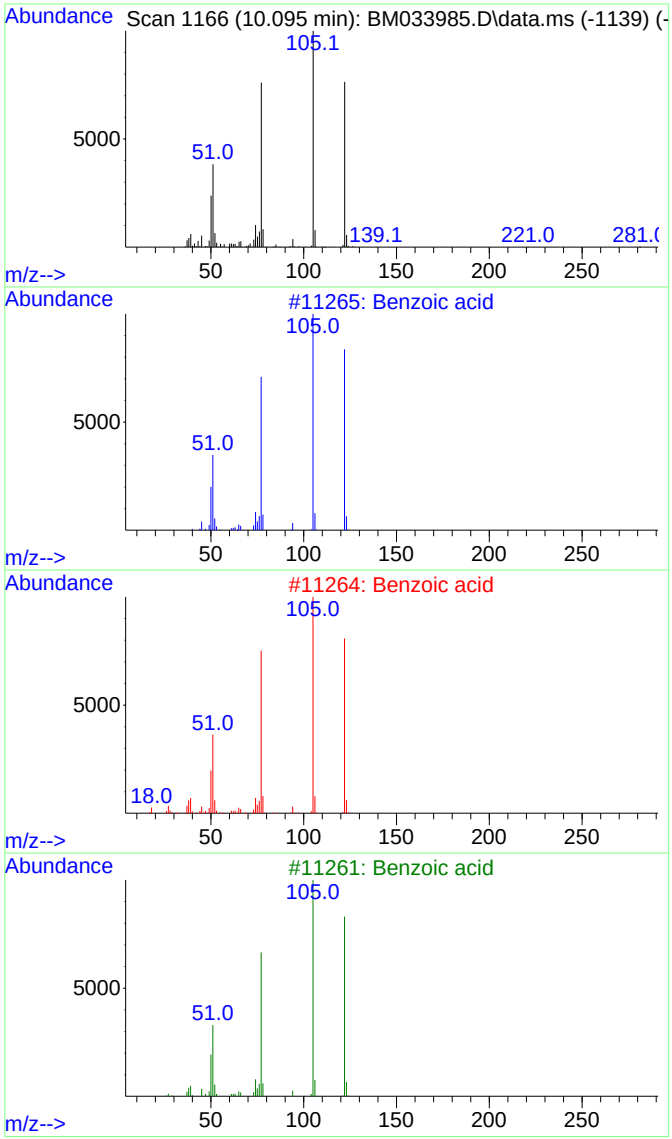
Quant Title : SVOA CALIBRATION

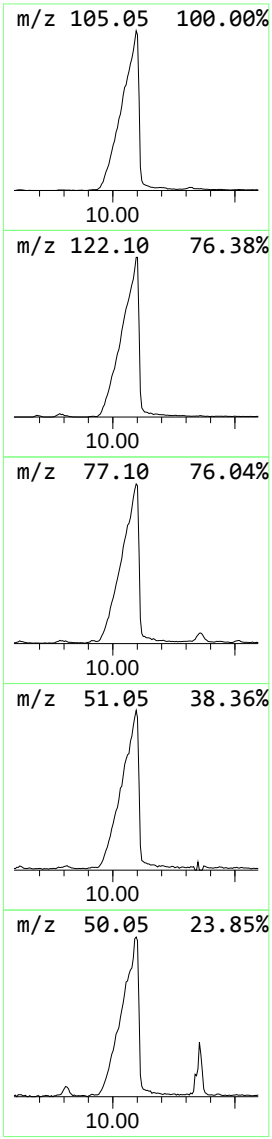
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 23 Benzoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.095	40.97 ng/ul	2721880	Naphthalene-d8	10.736

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzoic acid	122	C7H6O2	000065-85-0	96
2	Benzoic acid	122	C7H6O2	000065-85-0	94
3	Benzoic acid	122	C7H6O2	000065-85-0	94
4	Heptanediamide, N,N'-di-benzoyloxy-	398	C21H22N2O6	1000253-26-4	91
5	Benzoic acid	122	C7H6O2	000065-85-0	90





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ALS Vial : 56 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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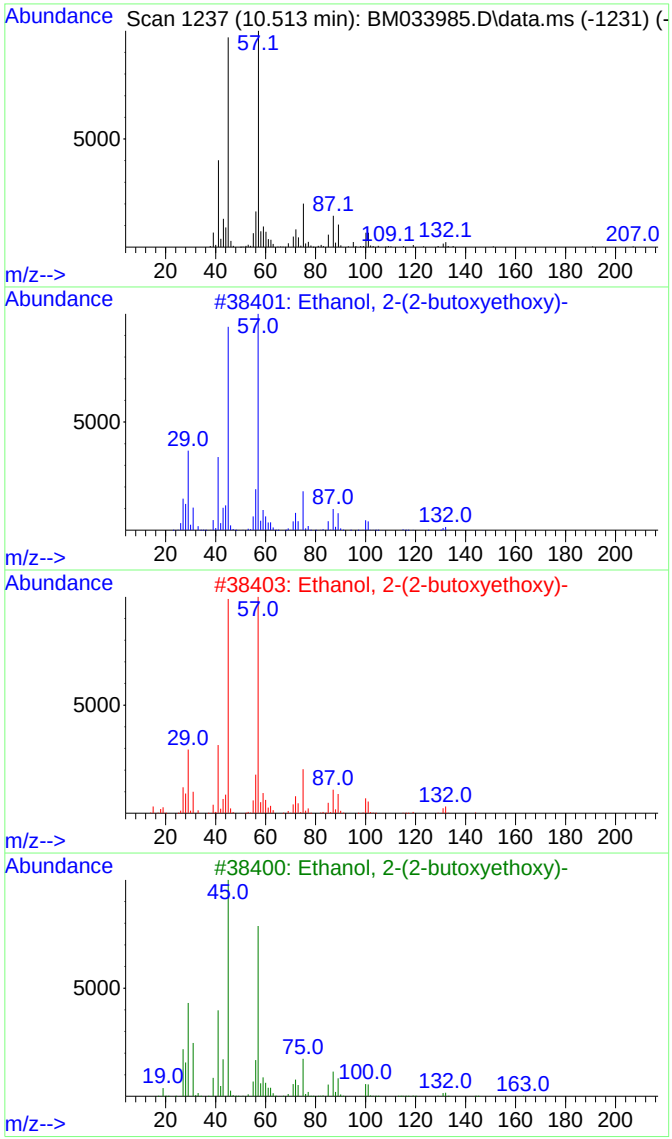
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM013122.M
Quant Title : SVOA CALIBRATION

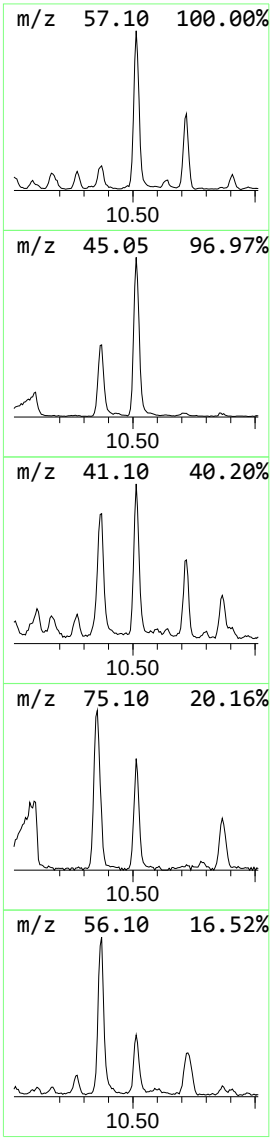
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 24 Ethanol, 2-(2-butoxyethoxy)- Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.513	5.51 ng/ul	366013	Naphthalene-d8	10.736

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
2	Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
3	Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
4	Ethanol, 1-(2-butoxyethoxy)-	162	C8H18O3	054446-78-5	83
5	Ethanol, 1-(2-butoxyethoxy)-	162	C8H18O3	054446-78-5	83





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Misc :
ALS Vial : 56 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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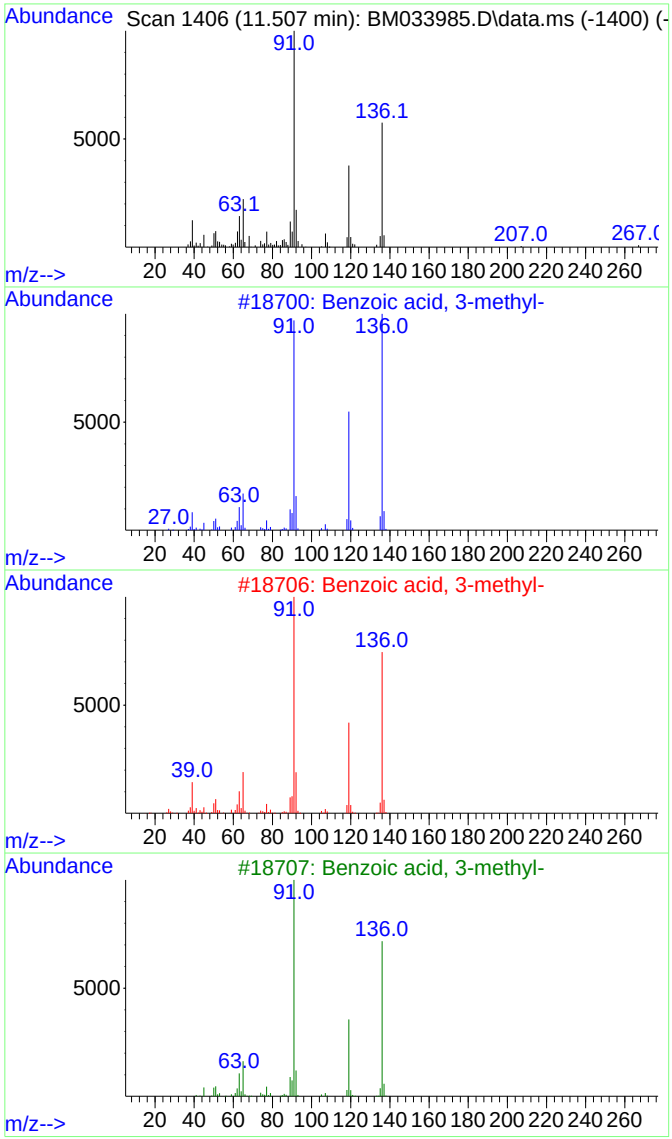
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM013122.M
Quant Title : SVOA CALIBRATION

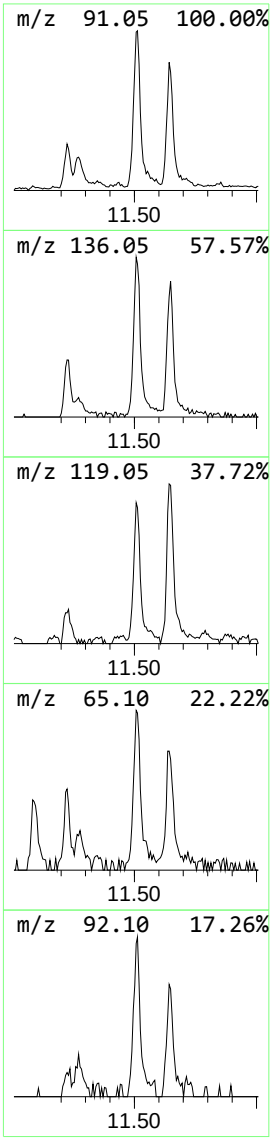
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 25 Benzoic acid, 3-methyl- Concentration Rank 31

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.507	2.47 ng/ul	164381	Naphthalene-d8	10.736

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzoic acid, 3-methyl-	136	C8H8O2	000099-04-7	94
2	Benzoic acid, 3-methyl-	136	C8H8O2	000099-04-7	93
3	Benzoic acid, 3-methyl-	136	C8H8O2	000099-04-7	91
4	Benzoic acid, 4-methyl-	136	C8H8O2	000099-94-5	91
5	Benzoic acid, 4-methyl-	136	C8H8O2	000099-94-5	91





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ALS Vial : 56 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

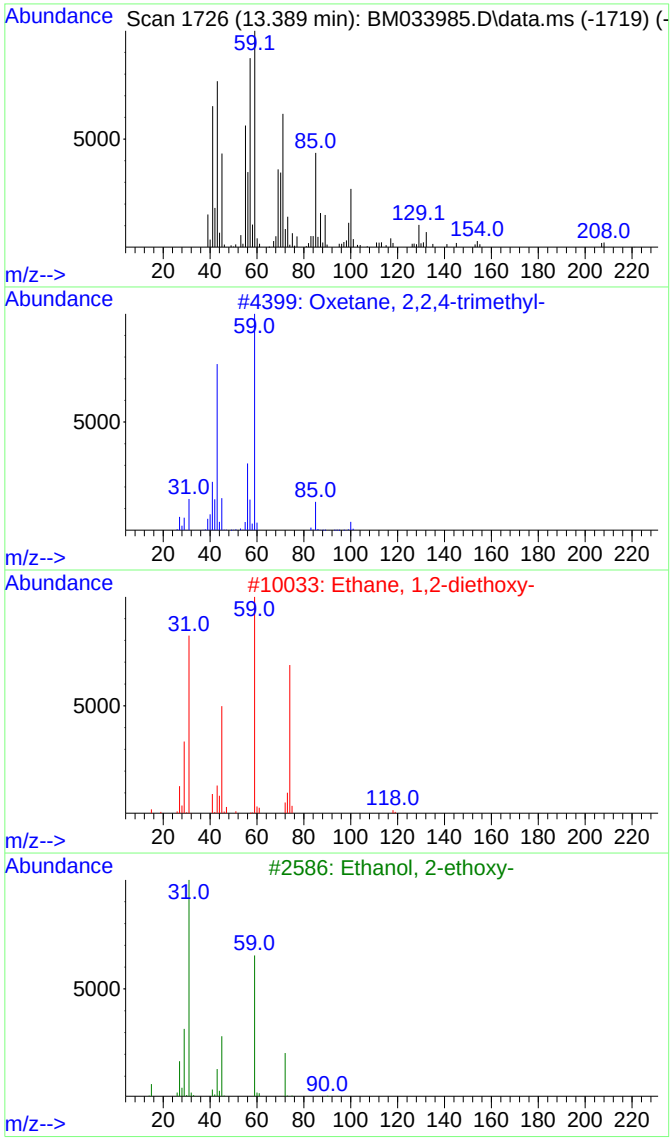
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

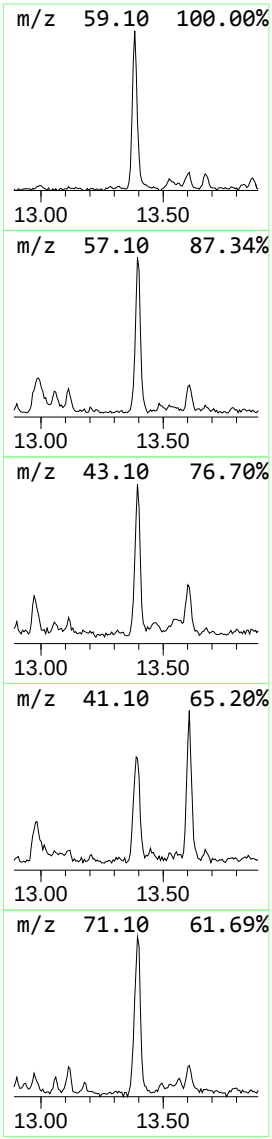
****

Peak Number 26 unknown-02 Concentration Rank 28

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.389	3.22 ng/ul	233733	Acenaphthene-d10	14.560

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Oxetane, 2,2,4-trimethyl-	100	C6H12O	023120-44-7	43
2	Ethane, 1,2-diethoxy-	118	C6H14O2	000629-14-1	38
3	Ethanol, 2-ethoxy-	90	C4H10O2	000110-80-5	35
4	Butanoic acid, 3-hydroxy-3-methy...	132	C6H12O3	006149-45-7	35
5	Oxalic acid, 6-ethyloct-3-yl pro...	272	C15H28O4	1000309-34-0	22





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ClientSampleId :
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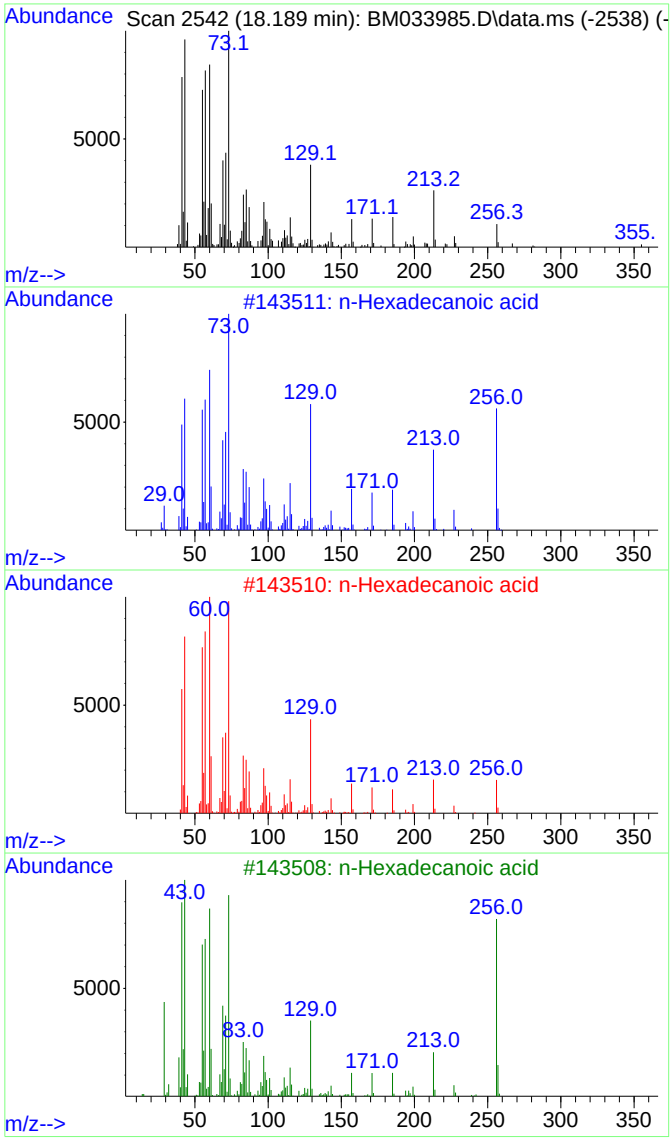
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM013122.M
Quant Title : SVOA CALIBRATION

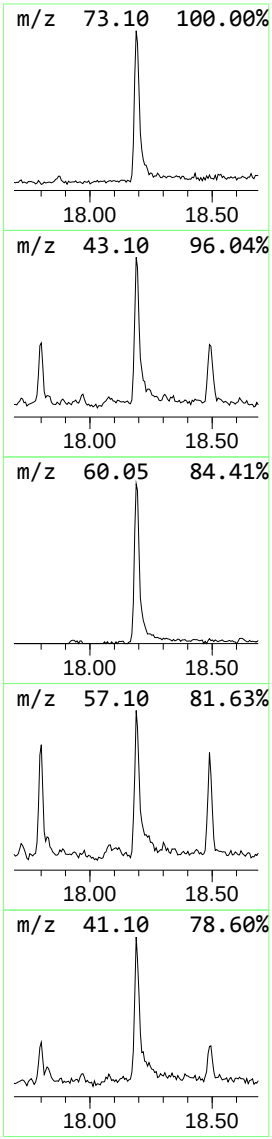
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 28 n-Hexadecanoic acid Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.189	5.31 ng/ul	383407	Phenanthrene-d10	17.295

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	Pentadecanoic acid	242	C15H30O2	001002-84-2	91





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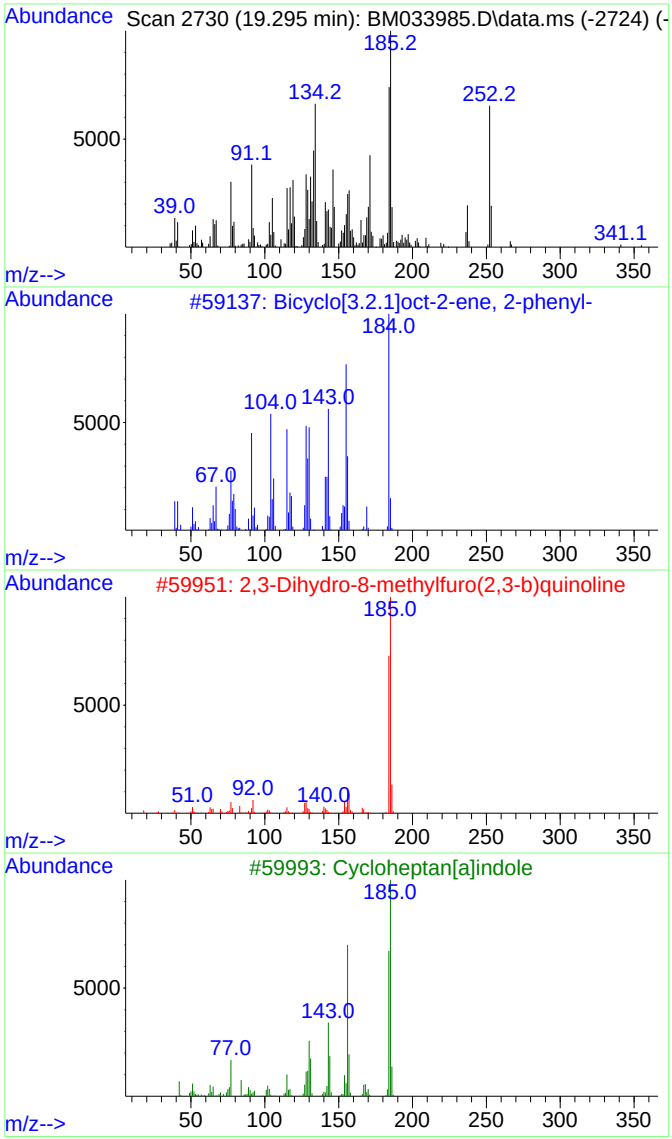
Quant Title : SVOA CALIBRATION

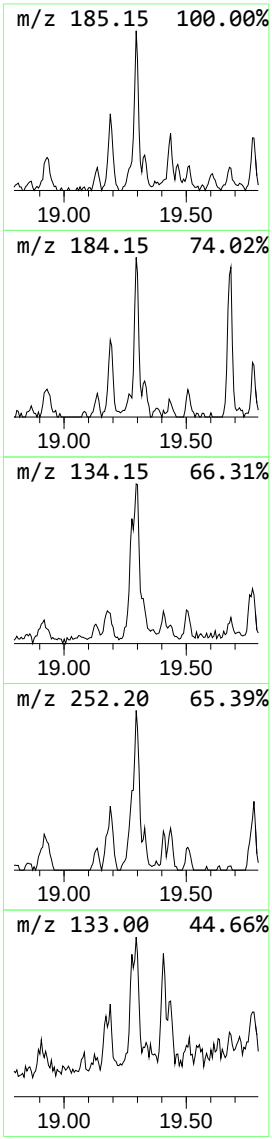
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 29 Bicyclo[3.2.1]oct-2-ene, 2-... Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.295	3.38 ng/ul	244127	Phenanthrene-d10	17.295

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Bicyclo[3.2.1]oct-2-ene, 2-phenyl-	184	C14H16	082027-19-8	53
2	2,3-Dihydro-8-methylfuro(2,3-b)q...	185	C12H11NO	064124-82-9	38
3	Cycloheptan[a]indole	185	C13H15N	002047-89-4	38
4	4-Pyridin-4-ylmethyl-phenol	185	C12H11NO	1000300-46-9	25
5	1,7-Trimethylene-2,3-dimethylindole	185	C13H15N	005825-43-4	18





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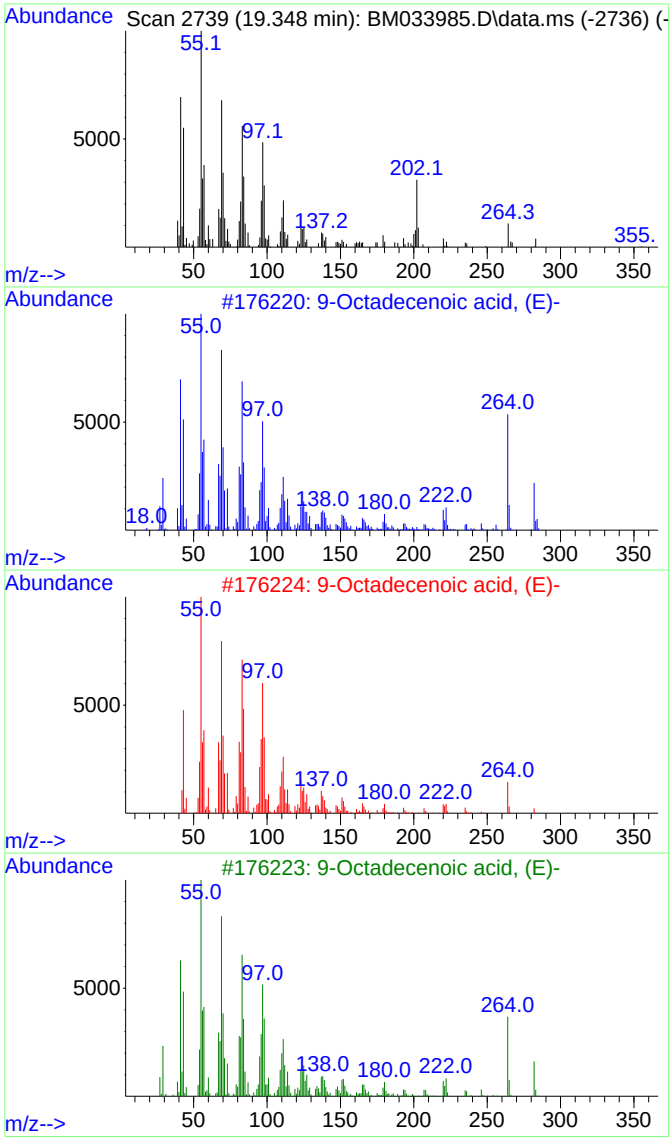
Quant Title : SVOA CALIBRATION

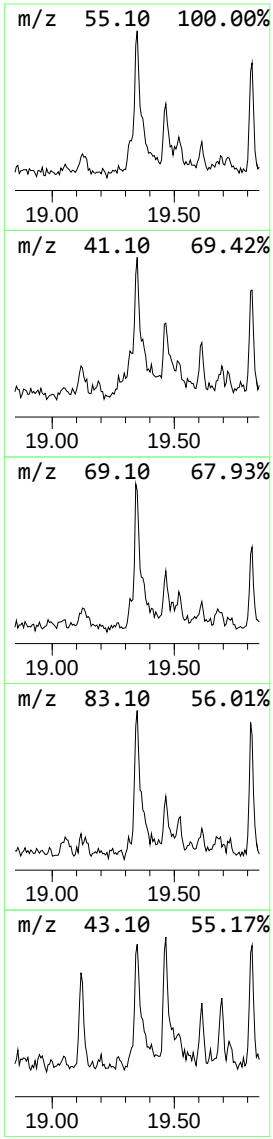
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 30 9-Octadecenoic acid, (E)- Concentration Rank 29

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.348	2.87 ng/ul	207468	Phenanthrene-d10	17.295

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9-Octadecenoic acid, (E)-	282	C18H34O2	000112-79-8	80
2	9-Octadecenoic acid, (E)-	282	C18H34O2	000112-79-8	74
3	9-Octadecenoic acid, (E)-	282	C18H34O2	000112-79-8	68
4	3-Bromobenzoic acid, hexadecyl e...	424	C23H37BrO2	070152-90-8	58
5	Cyclotetradecane	196	C14H28	000295-17-0	49





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ALS Vial : 56 Sample Multiplier: 1

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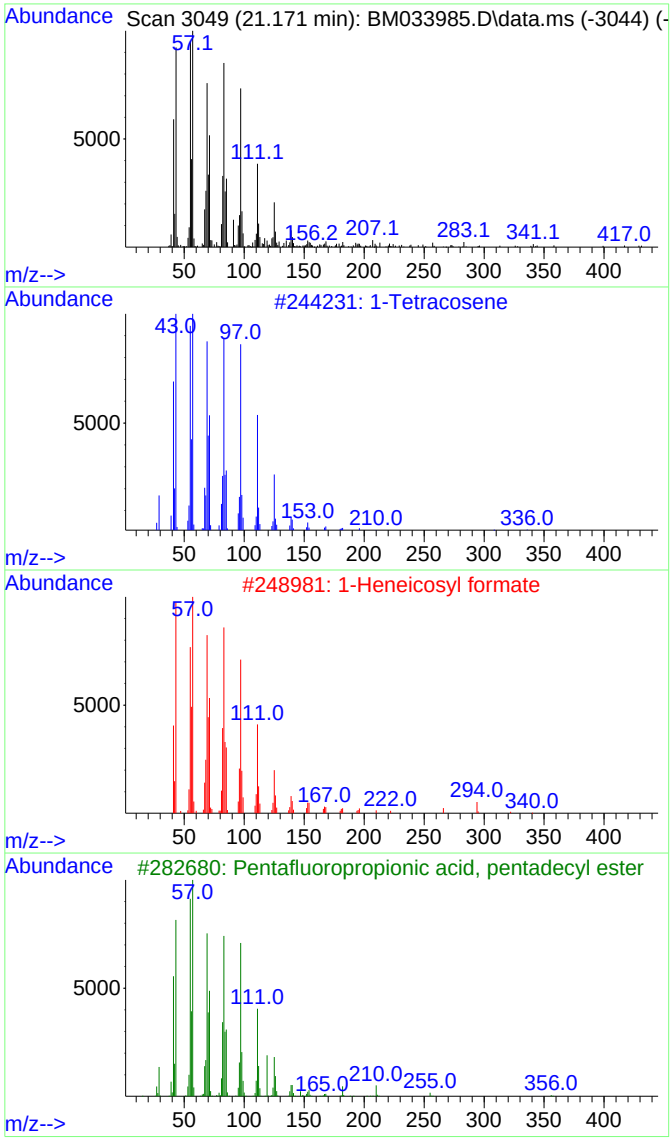
Quant Title : SVOA CALIBRATION

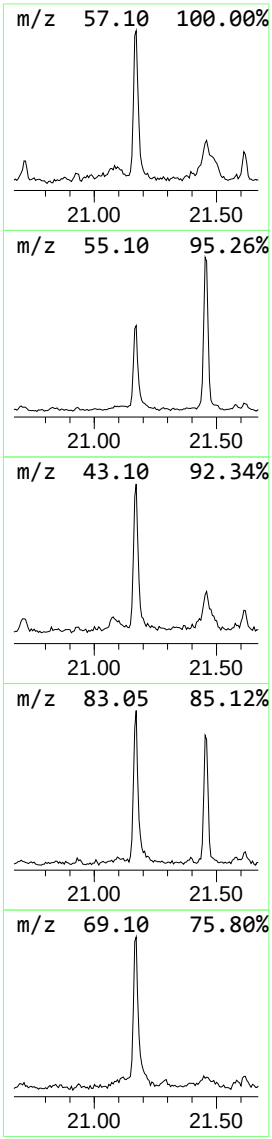
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 31 (DEL) Alkane: Straight-Chai... Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.171	4.16 ng/ul	443447	Chrysene-d12	21.459

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Tetracosene	336	C24H48	010192-32-2	95
2	1-Heneicosyl formate	340	C22H44O2	077899-03-7	94
3	Pentafluoropropionic acid, penta...	374	C18H31F5O2	959092-08-1	94
4	1-Hexacosene	364	C26H52	018835-33-1	94
5	Trifluoroacetoxy hexadecane	338	C18H33F3O2	006222-03-3	94





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Data File : BM033985.D
Acq On : 02 Feb 2022 21:59
Operator : CG/JU
Sample : N1355-09
Misc :
ALS Vial : 56 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0VF3

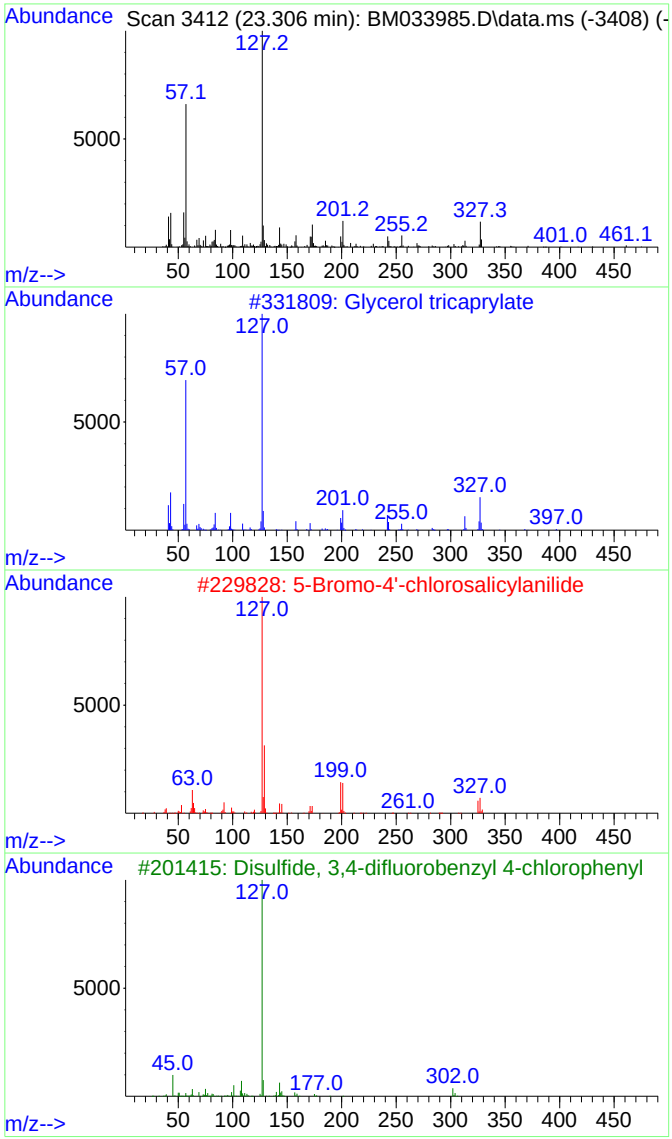
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM013122.M
Quant Title : SVOA CALIBRATION

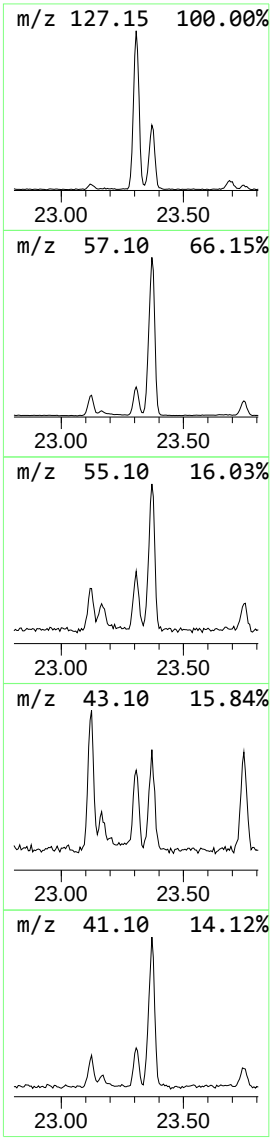
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 32 Glycerol tricaprylate Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.306	3.86 ng/ul	286649	Perylene-d12	23.847

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Glycerol tricaprylate	470	C27H50O6	000538-23-8	72
2	5-Bromo-4'-chlorosalicylanilide	325	C13H9BrClNO2	003679-64-9	64
3	Disulfide, 3,4-difluorobenzyl 4-...	302	C13H9ClF2S2	1000470-62-4	50
4	Fumaric acid, 4-bromophenyl ethy...	298	C12H11BrO4	1000344-75-3	50
5	Ethyl (2-amino-4-methyl-1,3-thia...	242	C10H14N2O3S	1000475-13-3	47





Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM020122\
Data File : BM033985.D
Acq On : 02 Feb 2022 21:59
Operator : CG/JU
Sample : N1355-09
Misc :
ALS Vial : 56 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0VF3

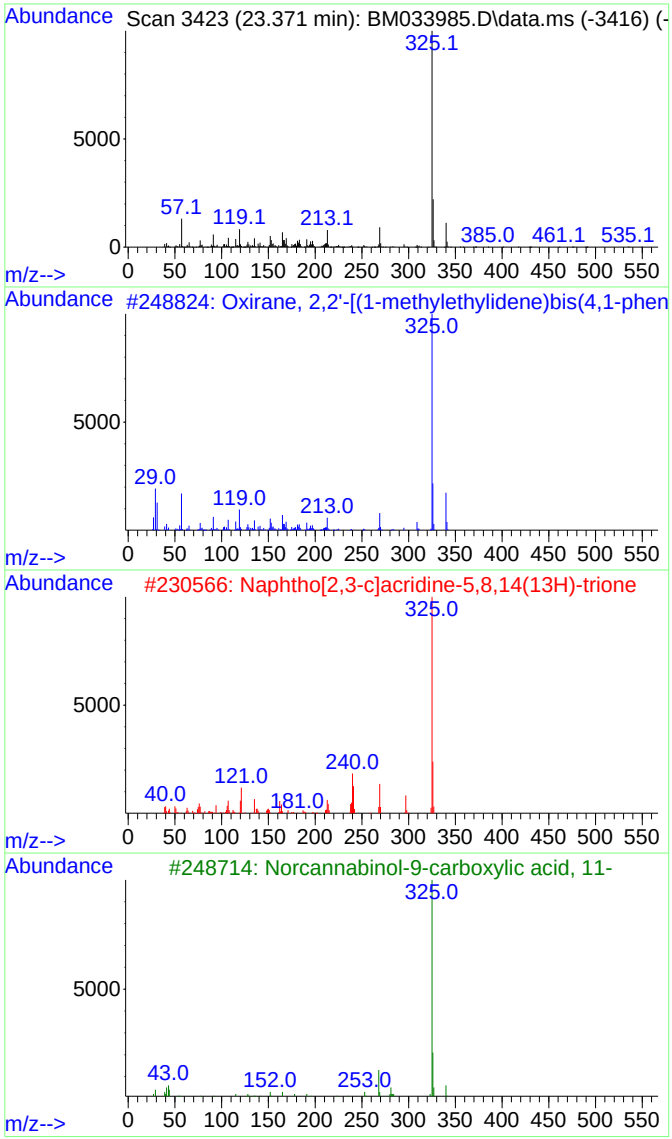
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM013122.M
Quant Title : SVOA CALIBRATION

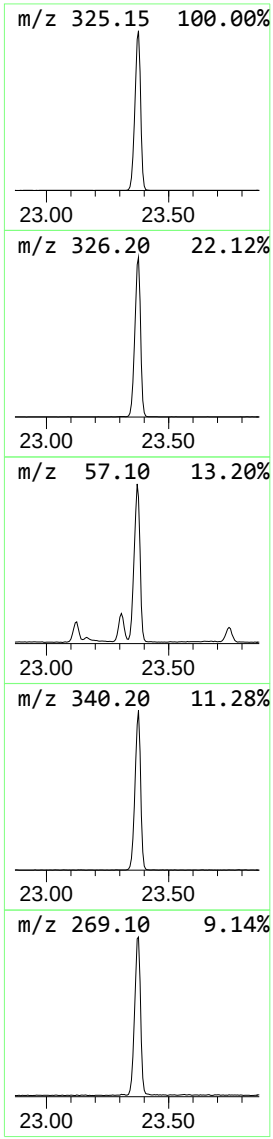
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 33 Oxirane, 2,2'-[(1-methyleth... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.371	97.73 ng/ul	7253290	Perylene-d12	23.847

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Oxirane, 2,2'-[(1-methylethylide...	340	C21H24O4	001675-54-3	90
2	Naphtho[2,3-c]acridine-5,8,14(13...	325	C21H11NO3	003569-01-5	72
3	Norcannabinol-9-carboxylic acid,...	340	C21H24O4	053989-32-5	64
4	2,6-Bis(1,1-dimethylethyl)cycloh...	325	C21H27NO2	017119-01-6	64
5	Bis(pentamethylcyclopentadienyl)...	325	C20H30Mn	067506-86-9	59





Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM020122\
Data File : BM033985.D
Acq On : 02 Feb 2022 21:59
Operator : CG/JU
Sample : N1355-09
Misc :
ALS Vial : 56 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0VF3

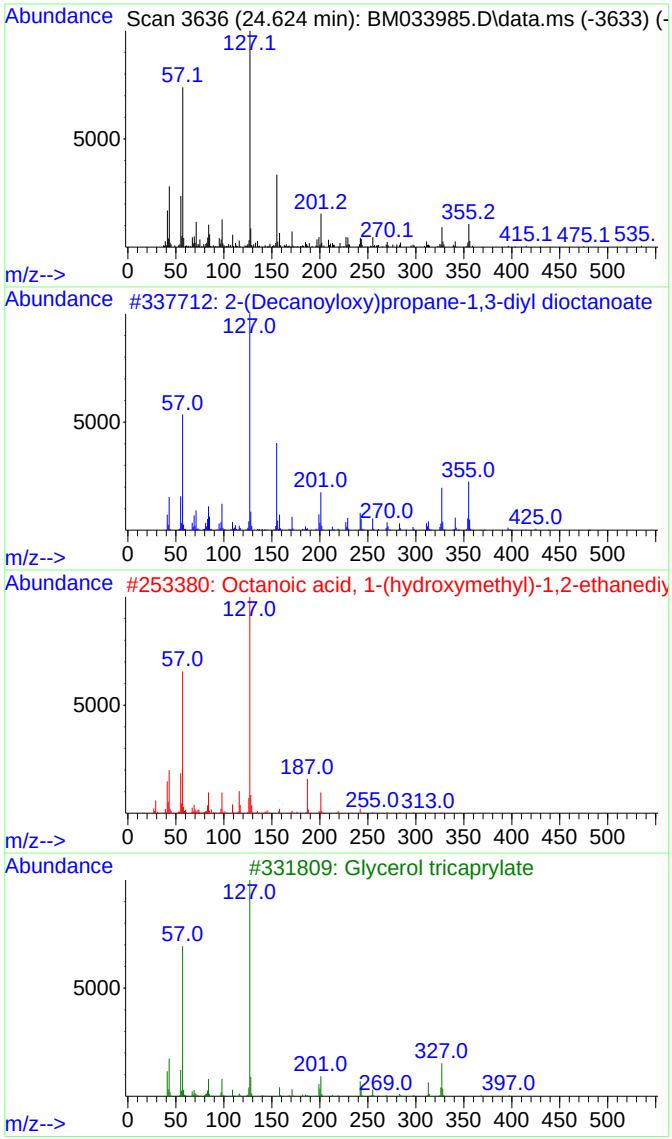
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM013122.M
Quant Title : SVOA CALIBRATION

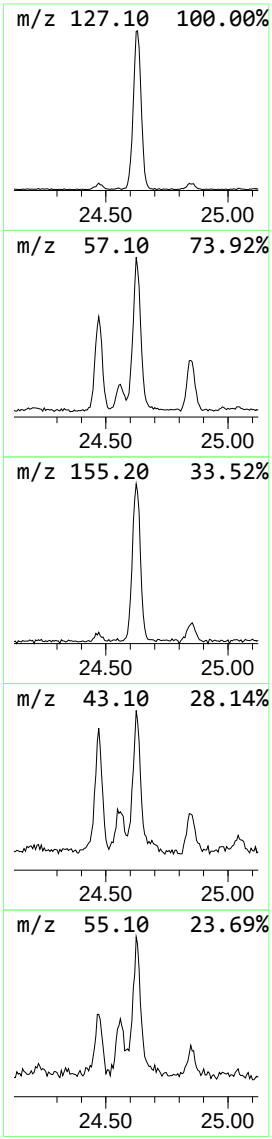
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 34 2-(Decanoyloxy)propane-1,3-... Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.624	6.14 ng/ul	455574	Perylene-d12	23.847

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-(Decanoyloxy)propane-1,3-diyl ...	498	C29H54O6	033368-87-5	72
2	Octanoic acid, 1-(hydroxymethyl)...	344	C19H36O5	060514-48-9	59
3	Glycerol tricaprylate	470	C27H50O6	000538-23-8	59
4	1,2-Dioctanoyl-sn-glycerol, O-ac...	386	C21H38O6	1000452-09-1	52
5	Octanoic acid, 2-formyl-4,6-dich...	316	C15H18Cl2O3	1000330-95-4	47





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Data File : BM033985.D
Acq On : 02 Feb 2022 21:59
Operator : CG/JU
Sample : N1355-09
Misc :
ALS Vial : 56 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0VF3

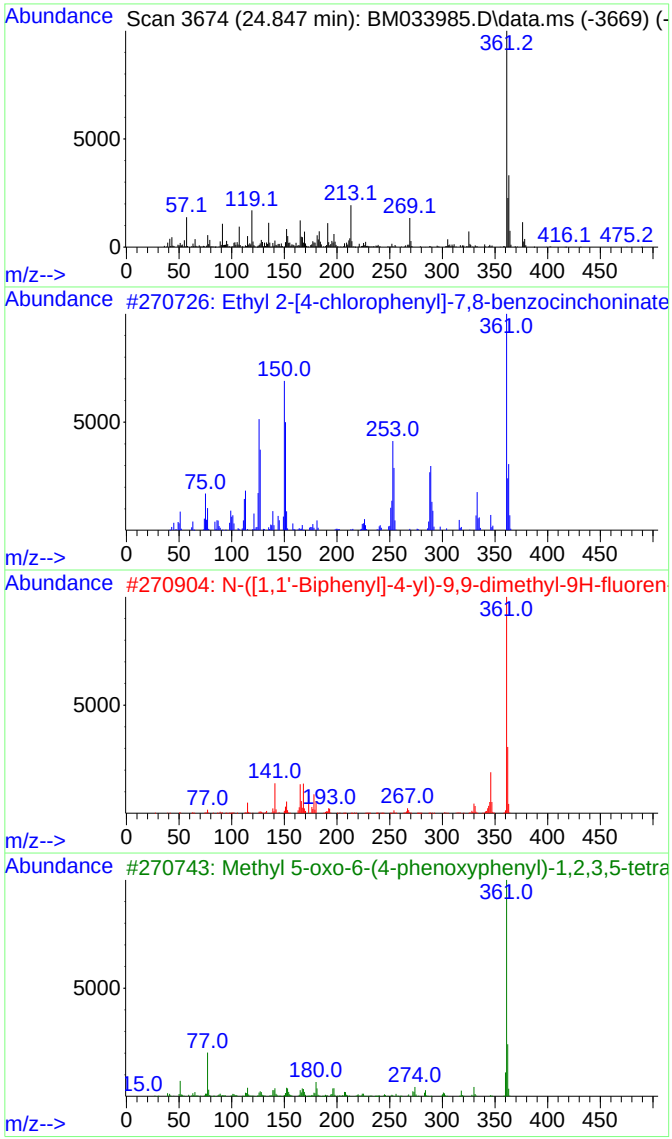
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM013122.M
Quant Title : SVOA CALIBRATION

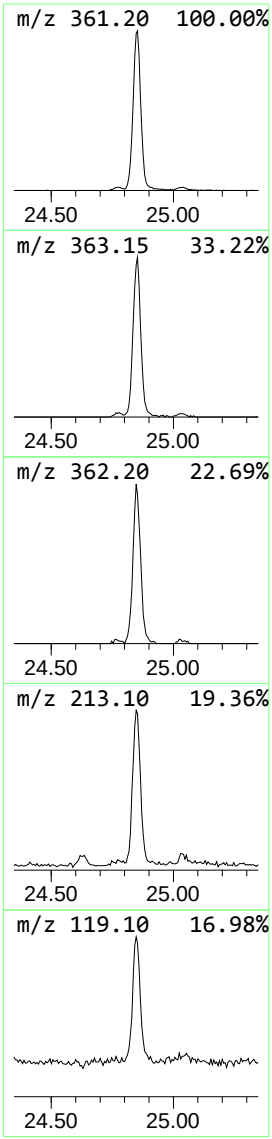
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 35 Ethyl 2-[4-chlorophenyl]-7,... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.847	12.01 ng/ul	891296	Perylene-d12	23.847

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethyl 2-[4-chlorophenyl]-7,8-ben...	361	C22H16ClNO2	1000256-80-8	93
2	N-([1,1'-Biphenyl]-4-yl)-9,9-dim...	361	C27H23N	897671-69-1	47
3	Methyl 5-oxo-6-(4-phenoxyphenyl)...	361	C22H19NO4	1000482-02-4	47
4	Maprotiline, N-chlorodifluoroace...	389	C22H22ClF2NO	1000448-25-9	42
5	Acifluorfen	361	C14H7ClF3NO5	050594-66-6	42





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM020122\
Data File : BM033985.D
Acq On : 02 Feb 2022 21:59
Operator : CG/JU
Sample : N1355-09
Misc :
ALS Vial : 56 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0VF3

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM013122.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
Propanoic acid,...	3.613	2.7	ng/ul	118922	1	7.925	867497	20.0
Methyl Isobutyl...	3.649	26.1	ng/ul	1130020	1	7.925	867497	20.0
2-Hexanol	3.872	3.8	ng/ul	166025	1	7.925	867497	20.0
Butanoic acid	4.060	30.5	ng/ul	1324600	1	7.925	867497	20.0
Ethanol, 2-prop...	4.472	36.4	ng/ul	1579170	1	7.925	867497	20.0
2-Pentanone, 4-...	5.013	67.5	ng/ul	2927880	1	7.925	867497	20.0
Ethylbenzene	5.407	6.0	ng/ul	262058	1	7.925	867497	20.0
Benzene, 1,3-di...	5.543	19.8	ng/ul	860322	1	7.925	867497	20.0
2-Heptanone	5.749	8.9	ng/ul	385866	1	7.925	867497	20.0
o-Xylene	5.919	12.6	ng/ul	548190	1	7.925	867497	20.0
Ethanol, 2-butoxy-	5.990	23.8	ng/ul	1031310	1	7.925	867497	20.0
2-Propanol, 1-b...	6.560	4.9	ng/ul	211179	1	7.925	867497	20.0
.beta.-Ethoxypr...	6.960	11.1	ng/ul	482211	1	7.925	867497	20.0
unknown-01	7.325	3.4	ng/ul	145406	1	7.925	867497	20.0
2-Propanol, 1-(...	7.525	44.0	ng/ul	1910180	1	7.925	867497	20.0
1-Propanol, 2-(...	7.719	11.8	ng/ul	513010	1	7.925	867497	20.0
1-Hexanol, 2-et...	7.995	2.3	ng/ul	101062	1	7.925	867497	20.0
Benzyl alcohol	8.190	310.3	ng/ul	13459200	1	7.925	867497	20.0
(DEL) Alkane: S...	8.466	2.3	ng/ul	101664	1	7.925	867497	20.0
(DEL) Alkane: S...	9.166	6.7	ng/ul	288284	1	7.925	867497	20.0
Hexanoic acid, ...	9.242	7.7	ng/ul	334702	1	7.925	867497	20.0
Benzoic acid	10.095	41.0	ng/ul	2721880	2	10.736	1328710	20.0
Ethanol, 2-(2-b...	10.513	5.5	ng/ul	366013	2	10.736	1328710	20.0
Benzoic acid, 3...	11.507	2.5	ng/ul	164381	2	10.736	1328710	20.0
unknown-02	13.389	3.2	ng/ul	233733	3	14.560	1450830	20.0
n-Hexadecanoic ...	18.189	5.3	ng/ul	383407	4	17.295	1445120	20.0
Bicyclo[3.2.1]o...	19.295	3.4	ng/ul	244127	4	17.295	1445120	20.0
9-Octadecenoic ...	19.348	2.9	ng/ul	207468	4	17.295	1445120	20.0
(DEL) Alkane: S...	21.171	4.2	ng/ul	443447	5	21.459	2130710	20.0
Glycerol tricap...	23.306	3.9	ng/ul	286649	6	23.847	1484420	20.0
Oxirane, 2,2'-[...	23.371	97.7	ng/ul	7253290	6	23.847	1484420	20.0
2-(Decanoyloxy)...	24.624	6.1	ng/ul	455574	6	23.847	1484420	20.0
Ethyl 2-[4-chlo...	24.847	12.0	ng/ul	891296	6	23.847	1484420	20.0