LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102121\
Data File : BP007554.D
Acq On : 21 Oct 2021 14:37
Operator : CG/JU
Sample : M4262-03
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : BFY56

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-BP101421.M
Title : SVOA CALIBRATION

Signal : TIC: BP007554.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.352	7	11	12	rBV4	7210	8383	0.16%	0.018%
2	3.375	12	15	18	rVV	43750	61972	1.19%	0.132%
3	3.411	18	21	32	rVB	256198	337838	6.48%	0.718%
4	3.664	59	64	72	rBV	40047	56985	1.09%	0.121%
5	3.811	87	89	97	rBV	38575	64349	1.23%	0.137%
6	3.999	116	121	125	rBV	42217	66293	1.27%	0.141%
7	4.040	125	128	133	rVB	27627	36361	0.70%	0.077%
8	4.128	140	143	147	rVB2	11916	14367	0.28%	0.031%
9	4.222	156	159	165	rVB4	10858	15457	0.30%	0.033%
10	4.799	249	257	263	rVV	62466	107580	2.06%	0.229%
11	4.952	272	283	290	rVB	77450	140385	2.69%	0.299%
12	5.052	296	300	306	rVB3	6957	12280	0.24%	0.026%
13	5.393	350	358	365	rBV	53170	95277	1.83%	0.203%
14	6.034	463	467	473	rVB3	10881	17167	0.33%	0.037%
15	6.340	510	519	532	rBV	88495	167216	3.21%	0.356%
16	6.952	615	623	628	rBV2	62616	108185	2.07%	0.230%
17	7.111	644	650	660	rVB	195944	325170	6.23%	0.691%
18	7.275	672	678	694	rVB	564318	897936	17.22%	1.910%
19	7.481	706	713	720	rVV	600703	971053	18.62%	2.065%
20	7.552	720	725	733	rVB	96318	147477	2.83%	0.314%
21	7.740	751	757	766	rVB	162958	246137	4.72%	0.523%
22	7.869	774	779	786	rBV2	15127	25864	0.50%	0.055%
23	7.952	786	793	800	rVV	591530	943495	18.09%	2.006%
24	8.022	800	805	810	rVB	18189	28110	0.54%	0.060%
25	8.199	830	835	842	rVB5	9467	16577	0.32%	0.035%
26	8.381	861	866	875	rVB7	6188	13803	0.26%	0.029%
27	8.510	885	888	892	rBV5	9264	14971	0.29%	0.032%
28	8.658	905	913	920	rBV	496916	816691	15.66%	1.737%
29	8.710	920	922	928	rVB3	12279	17047	0.33%	0.036%
30	9.105	980	989	1000	rBV	675628	1104834	21.18%	2.350%
31	9.228	1006	1010	1014	rVB6	6495	10452	0.20%	0.022%
32	9.287	1014	1020	1025	rBV4	11912	17921	0.34%	0.038%
33	9.575	1057	1069	1075	rVB2	16707	33144	0.64%	0.070%
34	9.834	1104	1113	1120	rBV	503375	828207	15.88%	1.761%
35	10.063	1147	1152	1153	rBV5	5537	6734	0.13%	0.014%
36	10.093	1156	1157	1161	rVB4	7603	7735	0.15%	0.016%

LSC Area Percent Report

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Sample : M4262-03
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : BFY56

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-BP101421.M
Title : SVOA CALIBRATION

37	10.375	1196	1205	1215	rBV	806044	1372394	26.31%	2.918%
38	10.446	1215	1217	1225	rVB9	4866	10370	0.20%	0.022%
39	10.540	1225	1233	1245	rBV	473485	782509	15.00%	1.664%
40	10.757	1256	1270	1281	rBV	792677	1327710	25.46%	2.823%
41	10.887	1283	1292	1300	rBV	103127	184653	3.54%	0.393%
42	11.110	1326	1330	1332	rBV5	4487	5852	0.11%	0.012%
43	11.169	1336	1340	1347	rVB9	4722	10268	0.20%	0.022%
44	11.275	1354	1358	1365	rVB3	5518	10884	0.21%	0.023%
45	11.410	1371	1381	1385	rBV6	6990	14883	0.29%	0.032%
46	11.499	1390	1396	1400	rBV7	6524	11401	0.22%	0.024%
47	11.546	1400	1404	1409	rBV6	5511	9136	0.18%	0.019%
48	11.951	1469	1473	1477	rBV2	24799	39061	0.75%	0.083%
49	11.999	1477	1481	1482	rVV2	28493	40349	0.77%	0.086%
50	12.016	1482	1484	1488	rVV	34119	46695	0.90%	0.099%
51	12.063	1488	1492	1501	rVB3	33048	65947	1.26%	0.140%
52	12.199	1510	1515	1518	rBV7	8565	13994	0.27%	0.030%
53	12.240	1518	1522	1526	rVV3	11913	20922	0.40%	0.044%
54	12.287	1526	1530	1534	rVV4	5563	9553	0.18%	0.020%
55	12.340	1534	1539	1544	rVB3	15579	25883	0.50%	0.055%
56	12.793	1612	1616	1619	rBV6	5224	6511	0.12%	0.014%
57	12.957	1639	1644	1649	rBV3	8582	13282	0.25%	0.028%
58	13.204	1682	1686	1691	rVV2	12872	18294	0.35%	0.039%
59	13.434	1720	1725	1729	rBV6	7973	11596	0.22%	0.025%
60	13.498	1729	1736	1741	rVV5	10688	19842	0.38%	0.042%
61	13.557	1742	1746	1751	rVB7	7147	10231	0.20%	0.022%
62	13.710	1764	1772	1776	rBV5	7716	13140	0.25%	0.028%
63	13.998	1814	1821	1827	rBV	1520325	2184021	41.87%	4.644%
64	14.051	1827	1830	1835	rVB6	14467	21411	0.41%	0.046%
65	14.275	1858	1868	1877	rBV	1650386	2549800	48.89%	5.422%
66	14.475	1898	1902	1904	rBV5	4928	7808	0.15%	0.017%
67	14.504	1904	1907	1911	rVB2	7538	8905	0.17%	0.019%
68	14.587	1913	1921	1931	rBV	1248415	1884393	36.13%	4.007%
69	14.769	1943	1952	1958	rBV	198648	304038	5.83%	0.647%
70	15.004	1988	1992	1995	rBV6	11705	17616	0.34%	0.037%
71	15.140	2008	2015	2040	rBV	165400	585379	11.22%	1.245%
72	15.404	2056	2060	2065	rBV2	21645	32343	0.62%	0.069%
73	15.457	2065	2069	2072	rVB4	9973	11088	0.21%	0.024%
74	15.581	2083	2090	2099	rBV	2266657	3320671	63.66%	7.062%
75	15.687	2103	2108	2115	rBV	549889	793769	15.22%	1.688%
76	15.822	2126	2131	2135	rBV	27021	37806	0.72%	0.080%

LSC Area Percent Report

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Acq On : 21 Oct 2021 14:37
Operator : CG/JU
Sample : M4262-03
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : BFY56

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-BP101421.M
Title : SVOA CALIBRATION

77	15.951	2149	2153	2155	rBV4	5382	7485	0.14%	0.016%
78	16.022	2160	2165	2170	rBV9	13043	25171	0.48%	0.054%
79	16.269	2202	2207	2209	rBV6	6886	12144	0.23%	0.026%
80	16.316	2213	2215	2218	rVB4	6157	6367	0.12%	0.014%
81	16.381	2223	2226	2230	rVB6	10314	15939	0.31%	0.034%
82	16.486	2242	2244	2249	rVB6	7504	9855	0.19%	0.021%
83	16.592	2259	2262	2266	rBV5	8918	12248	0.23%	0.026%
84	16.745	2284	2288	2293	rVB7	12189	18947	0.36%	0.040%
85	17.092	2344	2347	2350	rBV2	18482	26504	0.51%	0.056%
86	17.339	2383	2389	2396	rBV	1537346	2174348	41.69%	4.624%
87	17.439	2399	2406	2415	rBV2	2831827	4075535	78.14%	8.667%
88	18.022	2501	2505	2510	rVB	90563	106387	2.04%	0.226%
89	18.233	2537	2541	2550	rBV	348167	490262	9.40%	1.043%
90	19.257	2713	2715	2722	rVB5	35955	45163	0.87%	0.096%
91	19.328	2724	2727	2729	rBV2	38301	44592	0.85%	0.095%
92	19.463	2746	2750	2756	rBV2	178346	251001	4.81%	0.534%
93	19.669	2780	2785	2791	rBV	3658205	4971706	95.32%	10.573%
94	21.145	3032	3036	3042	rBV	487485	638321	12.24%	1.357%
95	21.422	3078	3083	3089	rBV	1888088	2549995	48.89%	5.423%
96	23.716	3466	3473	3483	rVB2	2559404	5215892	100.00%	11.092%
97	23.874	3494	3500	3510	rVB2	1310434	2688254	51.54%	5.717%

Sum of corrected areas: 47023977

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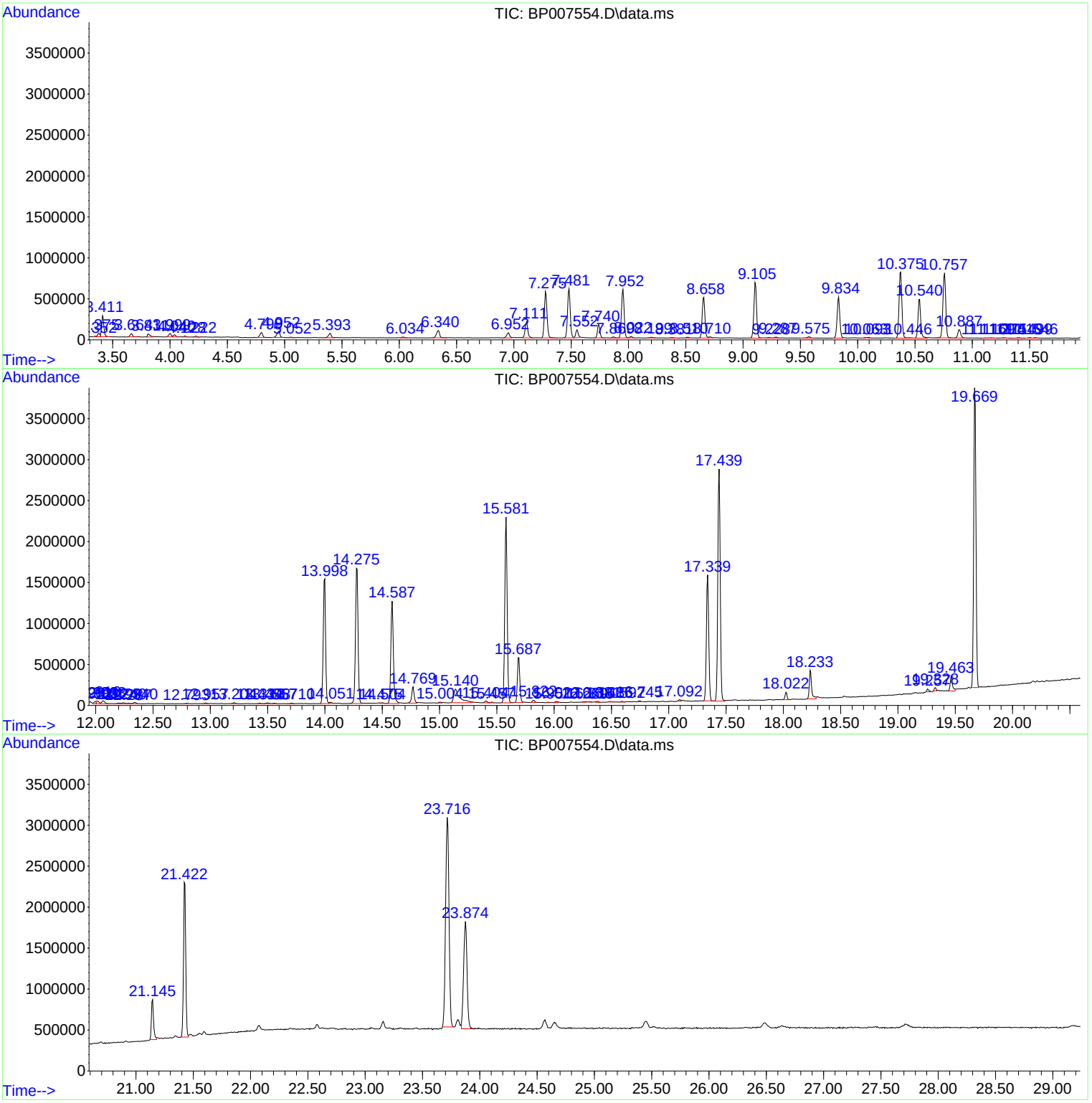
Data File : BP007554.D
Acq On : 21 Oct 2021 14:37
Operator : CG/JU
Sample : M4262-03
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BFY56

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP101421.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102121\
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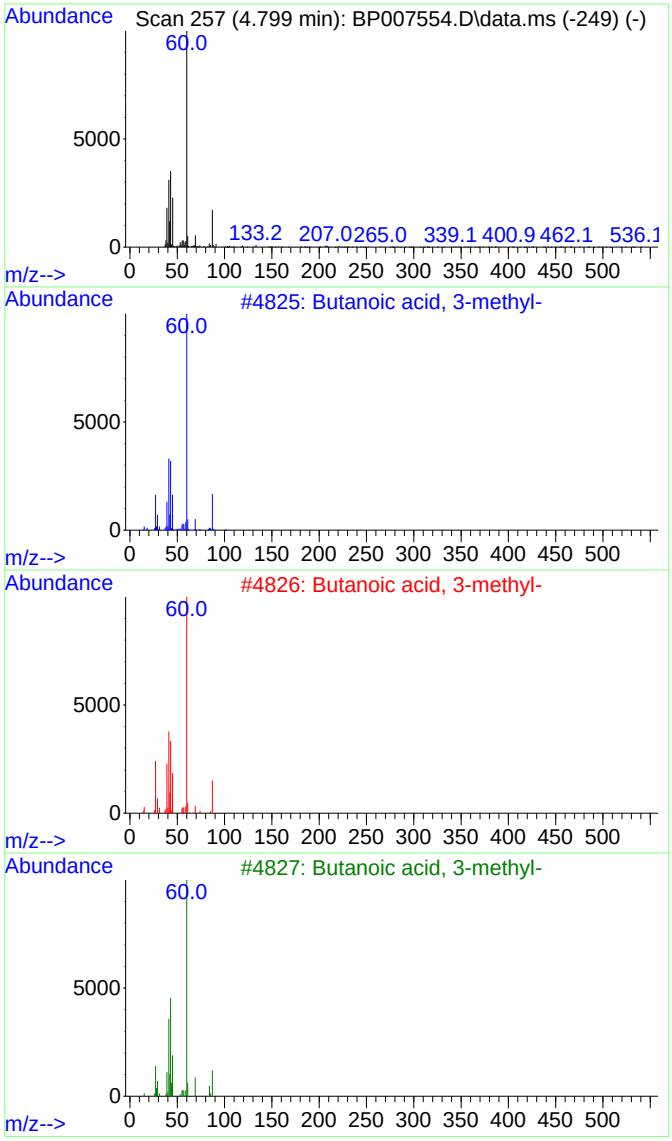
Quant Title : SVOA CALIBRATION

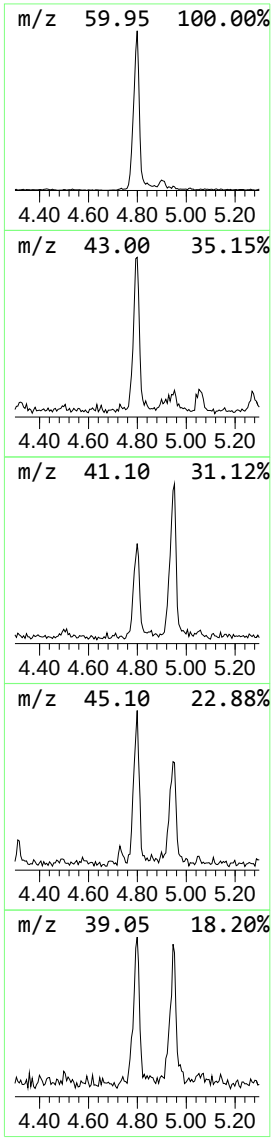
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 Butanoic acid, 3-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.799	2.28 ng/ul	107580	1,4-Dichlorobenzene-d4	7.952

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butanoic acid, 3-methyl-	102	C5H10O2	000503-74-2	64
2	Butanoic acid, 3-methyl-	102	C5H10O2	000503-74-2	64
3	Butanoic acid, 3-methyl-	102	C5H10O2	000503-74-2	50
4	Butanoic acid, 3-methyl-	102	C5H10O2	000503-74-2	43
5	Heptanoic acid	130	C7H14O2	000111-14-8	39





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Data File : BP007554.D
Acq On : 21 Oct 2021 14:37
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Sample : M4262-03
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BFY56

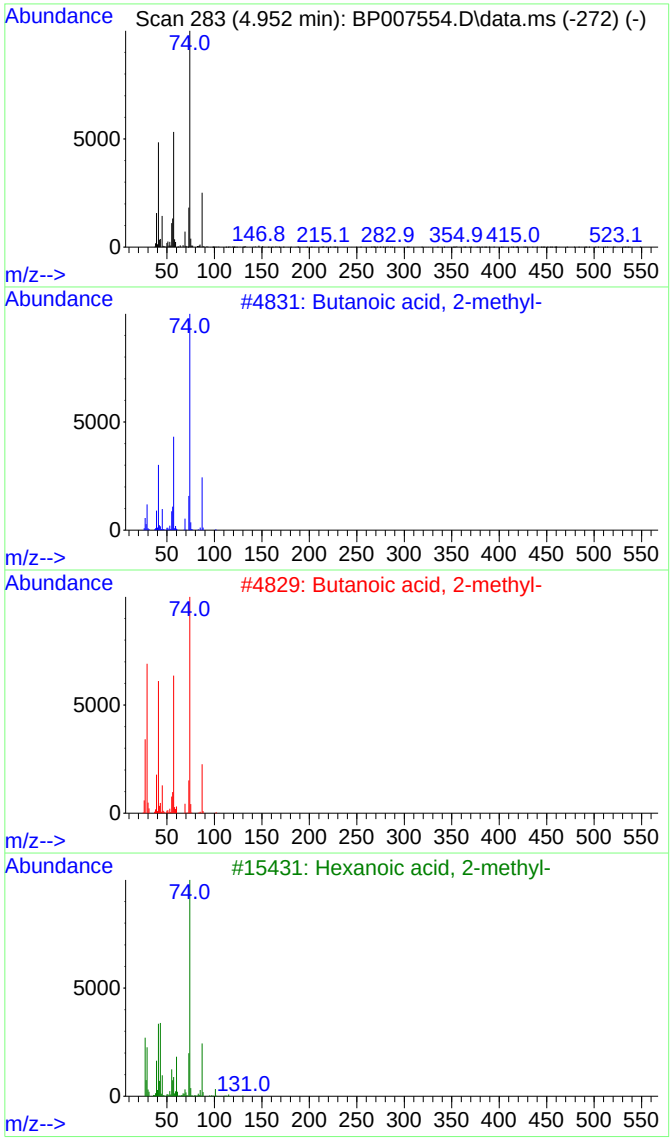
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP101421.M
Quant Title : SVOA CALIBRATION

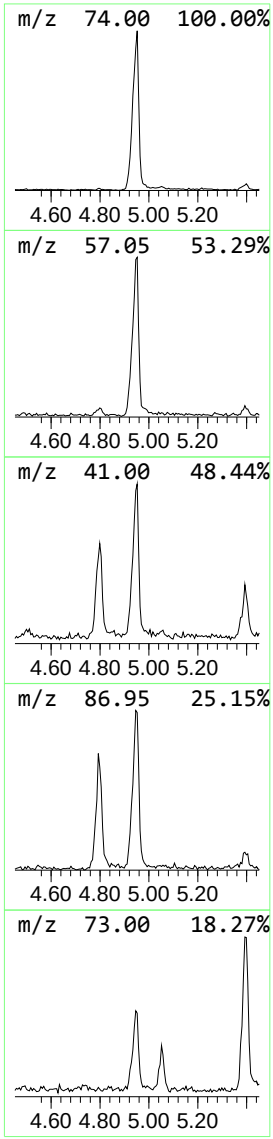
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 Butanoic acid, 2-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.952	2.98 ng/ul	140385	1,4-Dichlorobenzene-d4	7.952

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butanoic acid, 2-methyl-	102	C5H10O2	000116-53-0	90
2	Butanoic acid, 2-methyl-	102	C5H10O2	000116-53-0	86
3	Hexanoic acid, 2-methyl-	130	C7H14O2	004536-23-6	78
4	Butanoic acid, 2-methyl-	102	C5H10O2	000116-53-0	72
5	Butanoic acid, 2-methyl-	102	C5H10O2	000116-53-0	72





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ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP101421.M
Quant Title : SVOA CALIBRATION

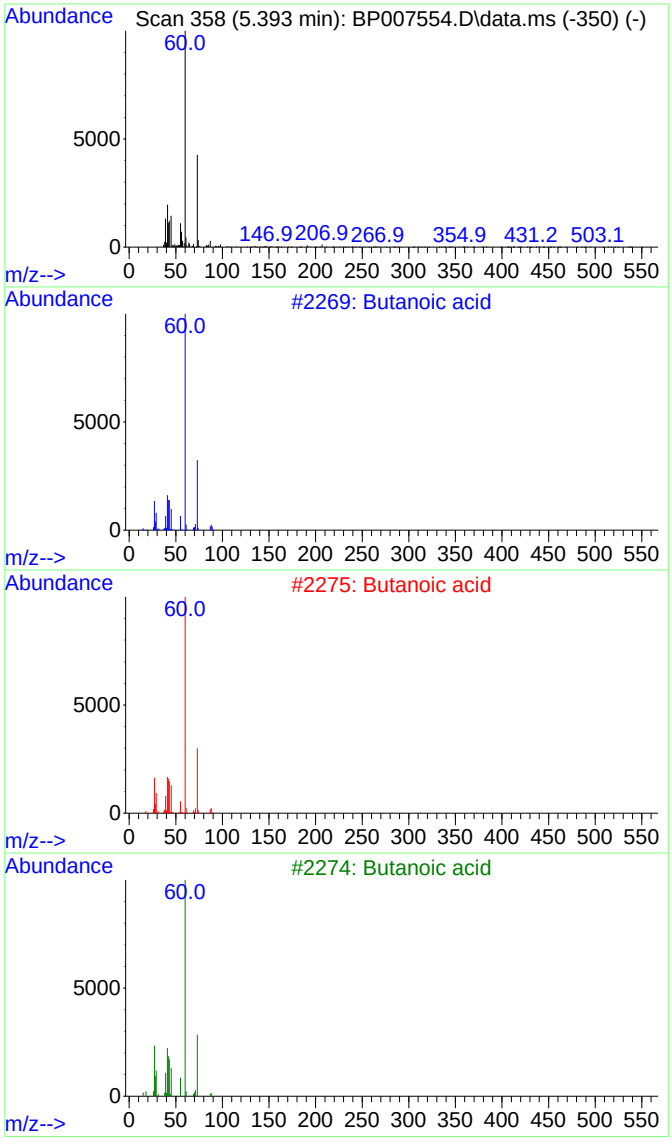
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

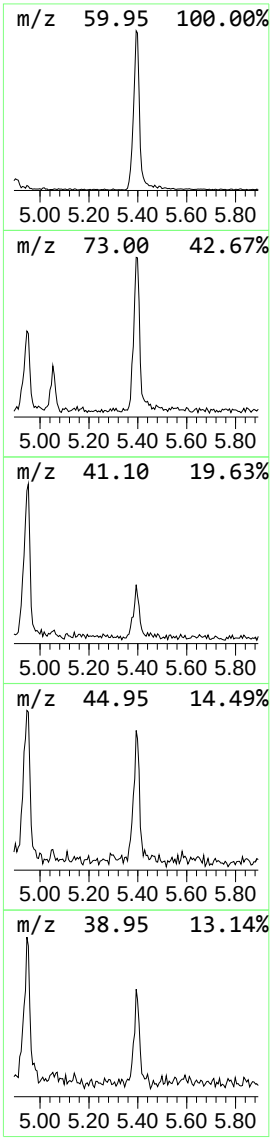
****

Peak Number 3 Butanoic acid Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.393	2.02 ng/ul	95277	1,4-Dichlorobenzene-d4	7.952

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butanoic acid	88	C4H8O2	000107-92-6	78
2	Butanoic acid	88	C4H8O2	000107-92-6	78
3	Butanoic acid	88	C4H8O2	000107-92-6	78
4	Pentanoic acid	102	C5H10O2	000109-52-4	78
5	Pentanoic acid	102	C5H10O2	000109-52-4	78





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ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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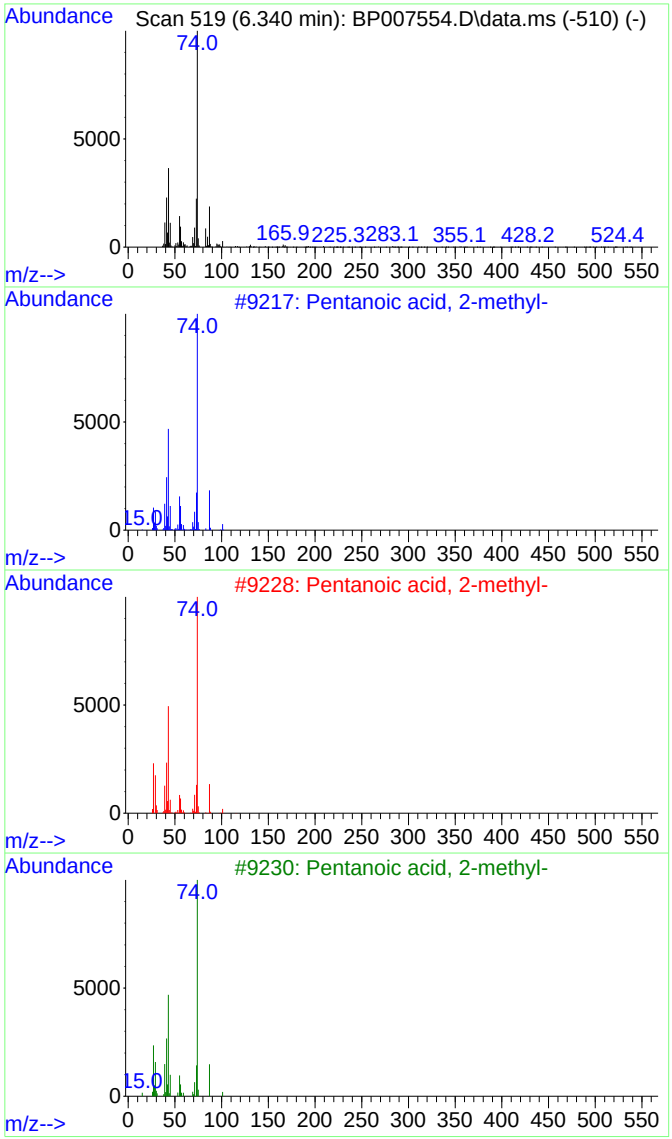
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP101421.M
Quant Title : SVOA CALIBRATION

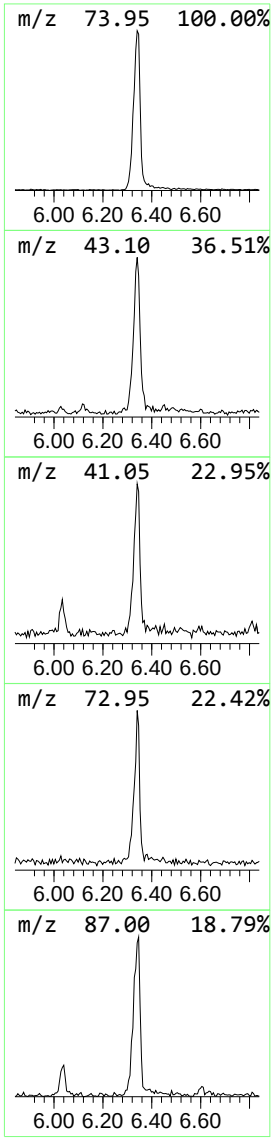
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Pentanoic acid, 2-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.340	3.54 ng/ul	167216	1,4-Dichlorobenzene-d4	7.952

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentanoic acid, 2-methyl-	116	C6H12O2	000097-61-0	78
2	Pentanoic acid, 2-methyl-	116	C6H12O2	000097-61-0	72
3	Pentanoic acid, 2-methyl-	116	C6H12O2	000097-61-0	64
4	Hexanoic acid, 2-methyl-	130	C7H14O2	004536-23-6	64
5	Pentanoic acid, 2-methyl-	116	C6H12O2	000097-61-0	64





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Acq On : 21 Oct 2021 14:37
Operator : CG/JU
Sample : M4262-03
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BFY56

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP101421.M
Quant Title : SVOA CALIBRATION

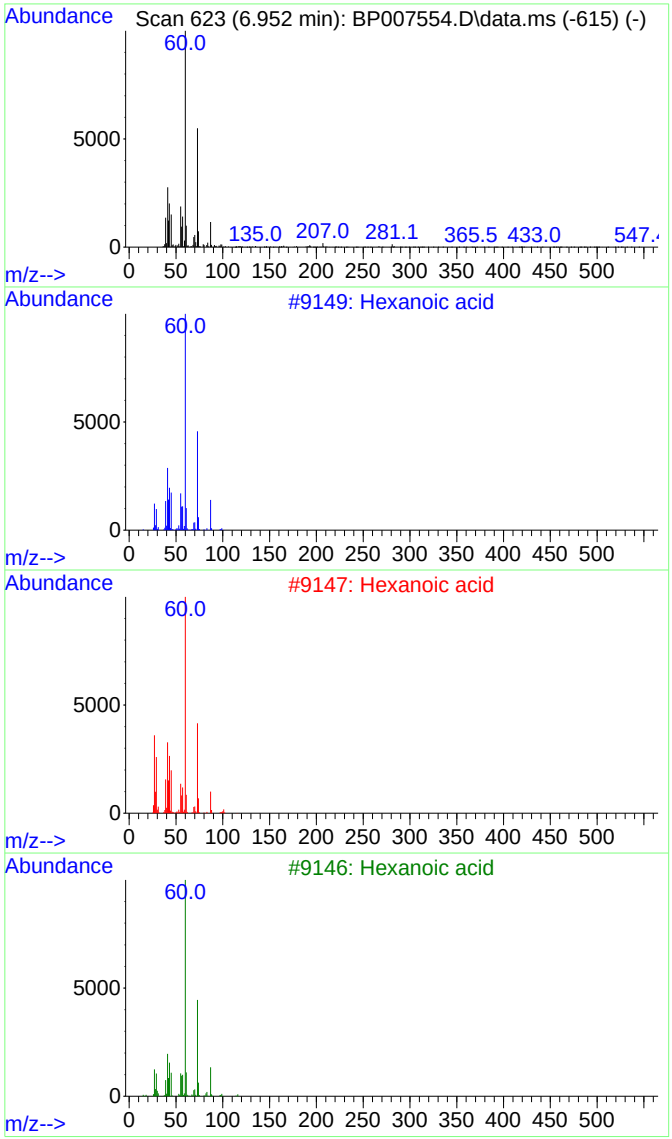
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

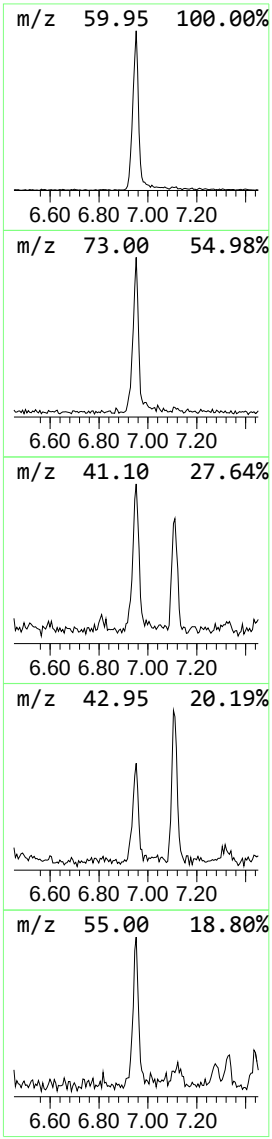
****

Peak Number 5 Hexanoic acid Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.952	2.29 ng/ul	108185	1,4-Dichlorobenzene-d4	7.952

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexanoic acid	116	C6H12O2	000142-62-1	90
2	Hexanoic acid	116	C6H12O2	000142-62-1	83
3	Hexanoic acid	116	C6H12O2	000142-62-1	83
4	Pentanoic acid	102	C5H10O2	000109-52-4	78
5	Octanoic acid	144	C8H16O2	000124-07-2	74





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Operator : CG/JU
Sample : M4262-03
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BFY56

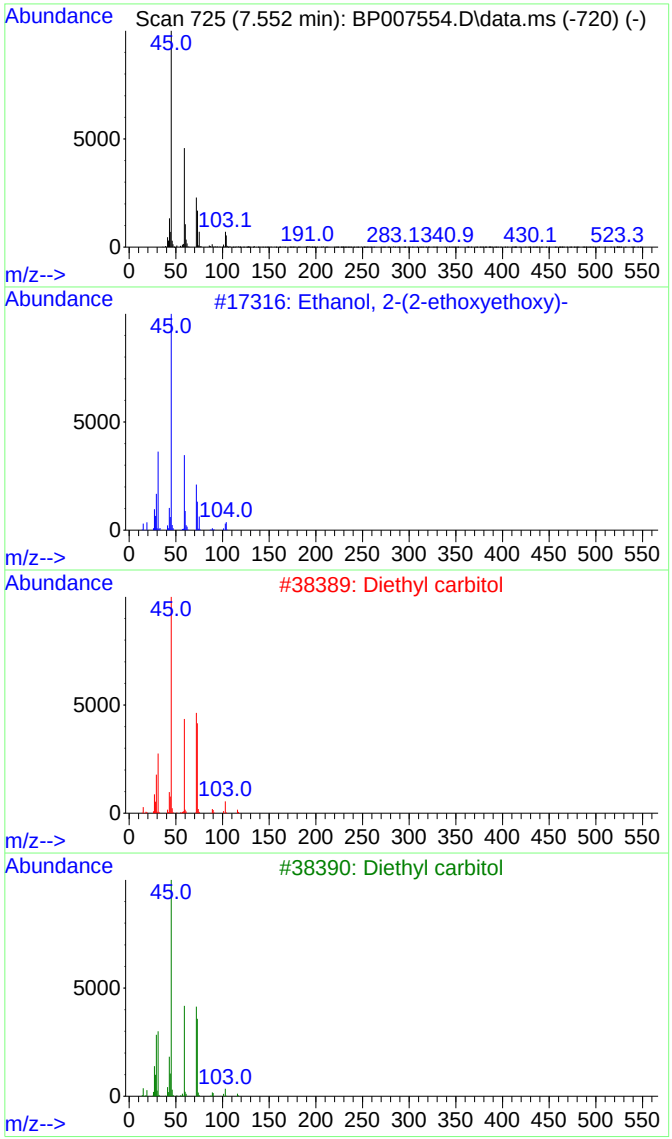
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP101421.M
Quant Title : SVOA CALIBRATION

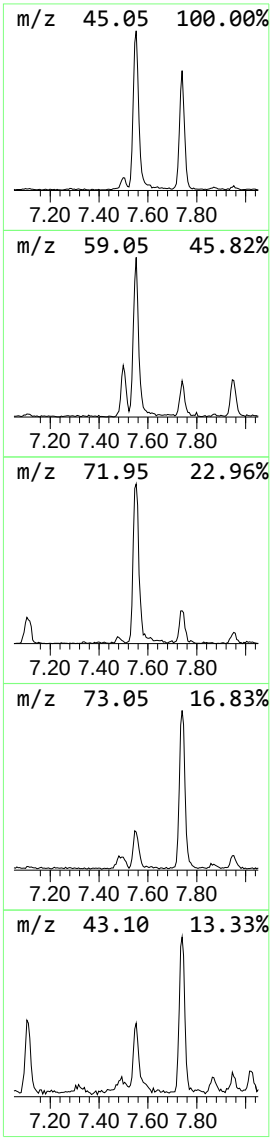
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 Ethanol, 2-(2-ethoxyethoxy)- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.552	3.13 ng/ul	147477	1,4-Dichlorobenzene-d4	7.952

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethanol, 2-(2-ethoxyethoxy)-	134	C6H14O3	000111-90-0	83
2	Diethyl carbitol	162	C8H18O3	000112-36-7	72
3	Diethyl carbitol	162	C8H18O3	000112-36-7	72
4	Ethanol, 2-(2-ethoxyethoxy)-	134	C6H14O3	000111-90-0	72
5	Ethanol, 2-(2-ethoxyethoxy)-	134	C6H14O3	000111-90-0	72





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Acq On : 21 Oct 2021 14:37
Operator : CG/JU
Sample : M4262-03
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BFY56

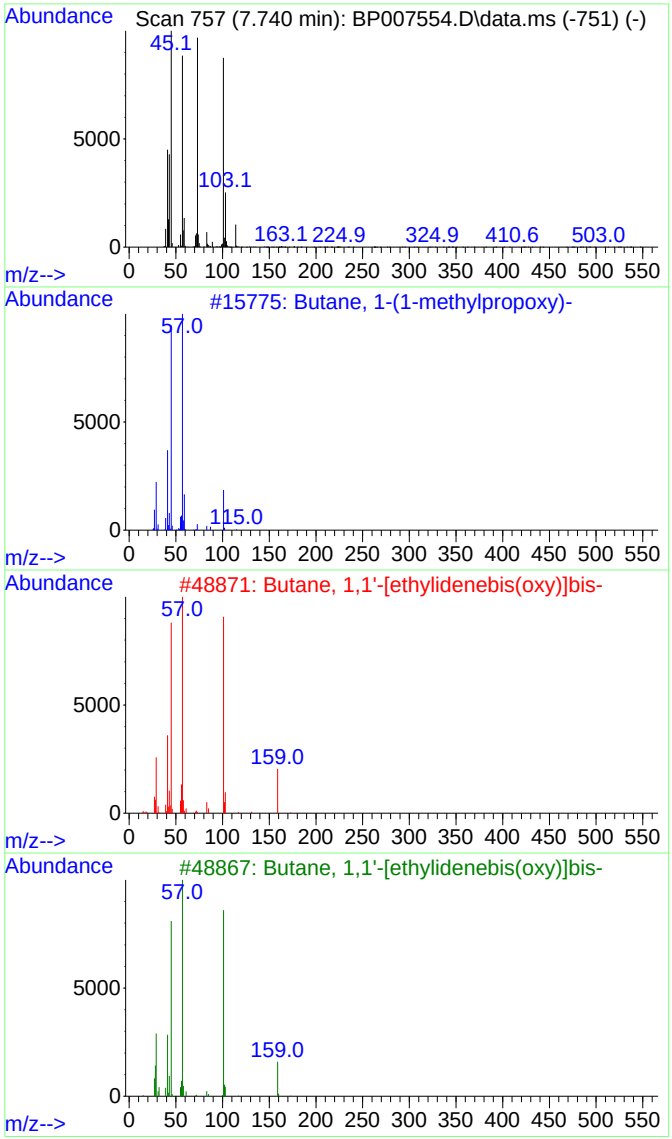
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP101421.M
Quant Title : SVOA CALIBRATION

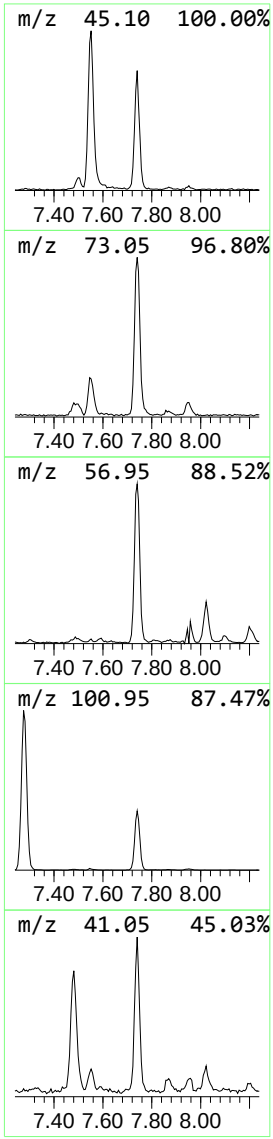
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 Butane, 1-(1-methylpropoxy)- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.740	5.22 ng/ul	246137	1,4-Dichlorobenzene-d4	7.952

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 1-(1-methylpropoxy)-	130	C8H18O	000999-65-5	64
2	Butane, 1,1'-[ethylidenebis(oxy)...	174	C10H22O2	000871-22-7	59
3	Butane, 1,1'-[ethylidenebis(oxy)...	174	C10H22O2	000871-22-7	59
4	2-t-Butyl-5-methyl[1,3]dioxolan-...	158	C8H14O3	130930-47-1	56
5	DL-Lactamide, butyl ether	145	C7H15NO2	1000452-56-5	50





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Sample : M4262-03
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : BFY56

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP101421.M
Quant Title : SVOA CALIBRATION

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

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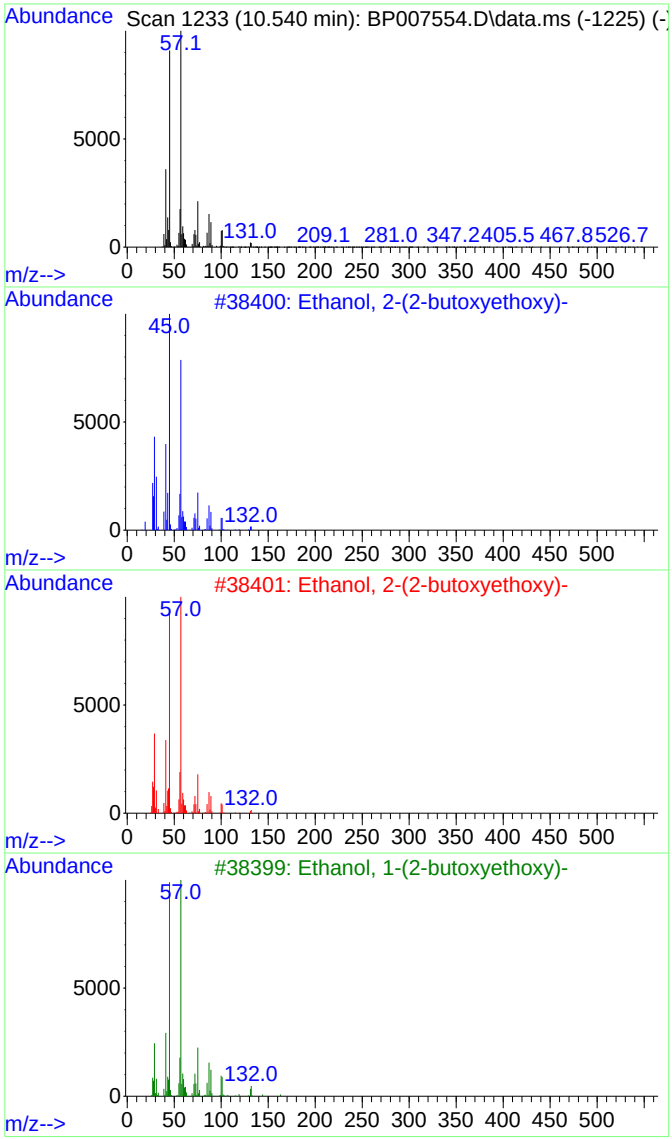
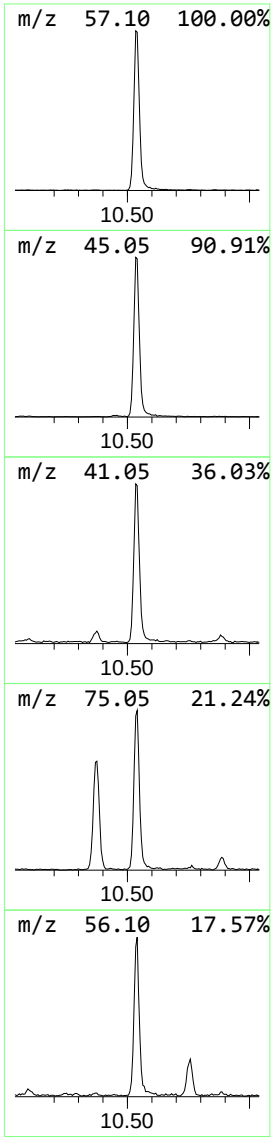
Peak Number 8 Ethanol, 2-(2-butoxyethoxy)- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.540	11.79 ng/ul	782509	Naphthalene-d8	10.757

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, 1-(2-butoxyethoxy)-	162	C8H18O3	054446-78-5	90
4	Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	83
5	Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	83





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Sample : M4262-03
Misc :
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Instrument :
BNA_P
ClientSampleId :
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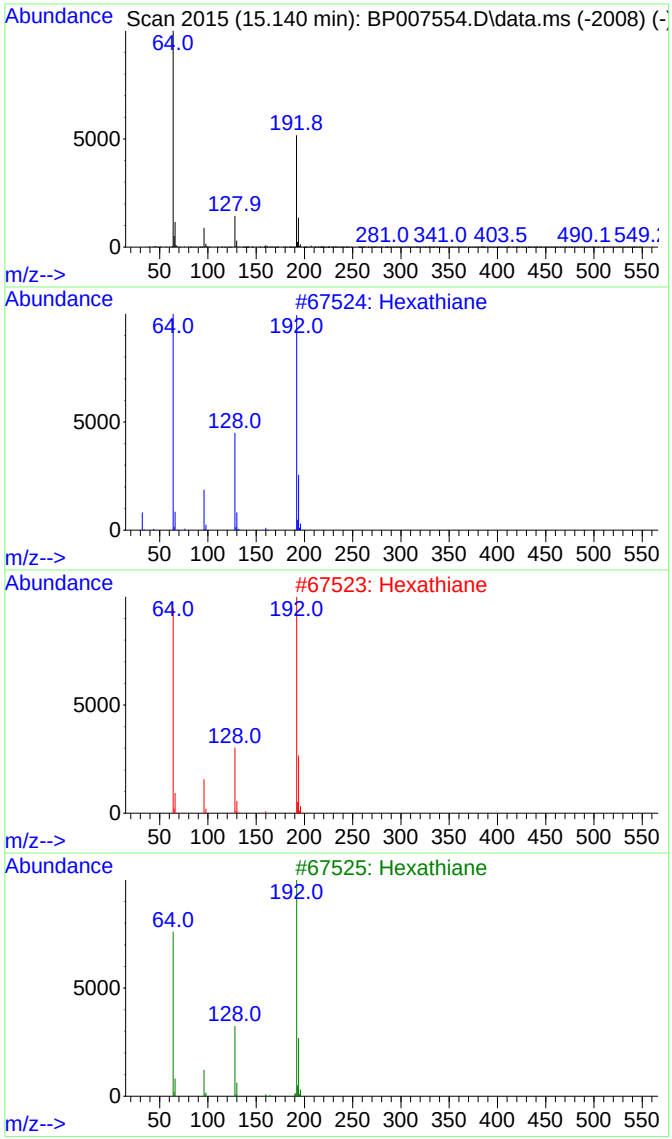
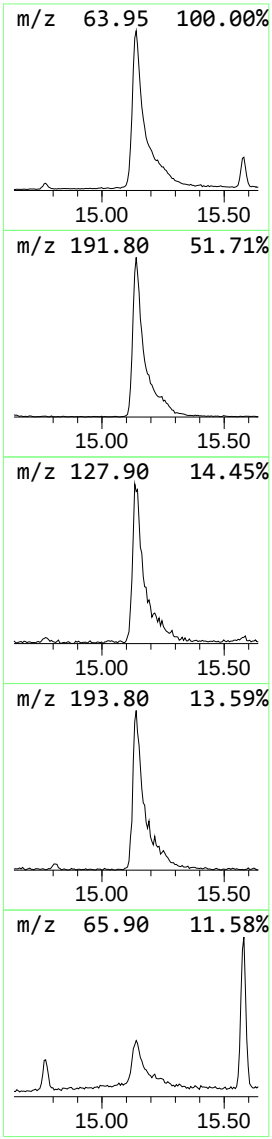
Peak Number 9 Hexathiane Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.140	6.21 ng/ul	585379	Acenaphthene-d10	14.587

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexathiane	192	S6	013798-23-7	91
2	Hexathiane	192	S6	013798-23-7	91
3	Hexathiane	192	S6	013798-23-7	91
4	Di(tert-butyl) trisulfide, perfl...	534	C8F18S3	016005-64-4	9
5	2-Chloro-7-methoxynaphthalene	192	C11H9ClO	067061-67-0	7





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ClientSampleId : BFY56

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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
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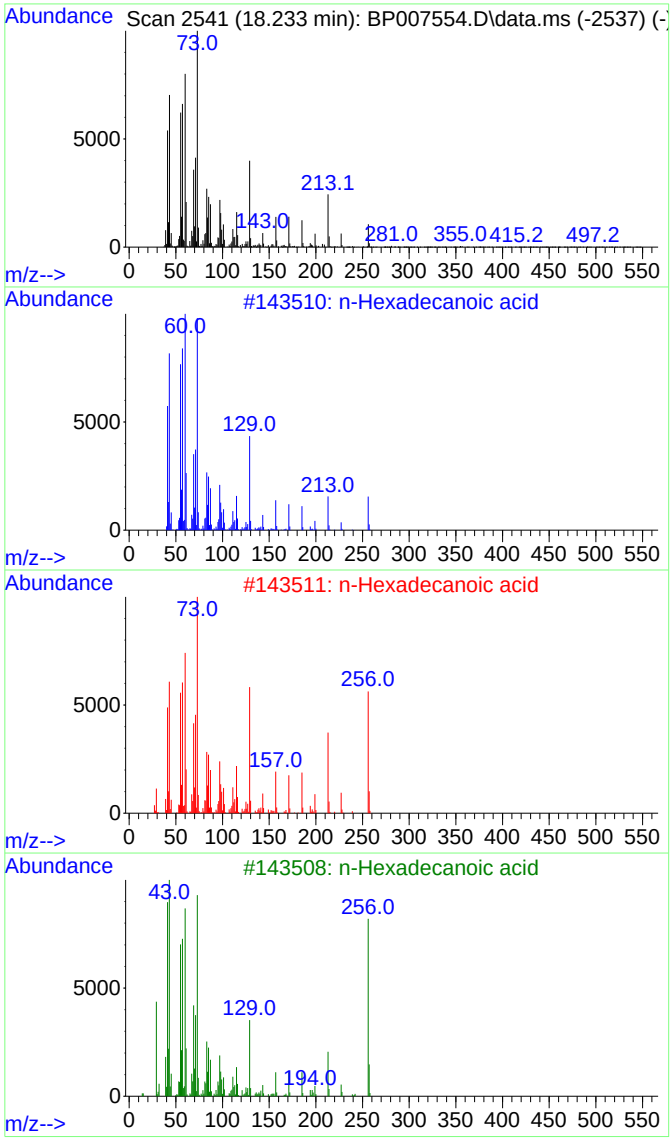
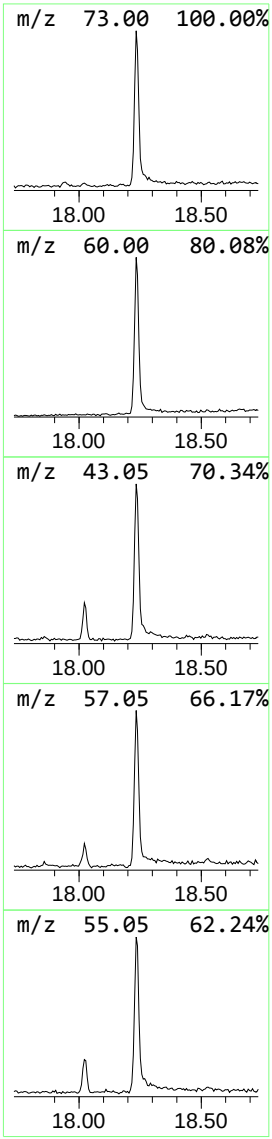
Peak Number 10 n-Hexadecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.233	4.51 ng/ul	490262	Phenanthrene-d10	17.339

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	Tridecanoic acid	214	C13H26O2	000638-53-9	93





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Sample : M4262-03
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : BFY56

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Quant Title : SVOA CALIBRATION

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

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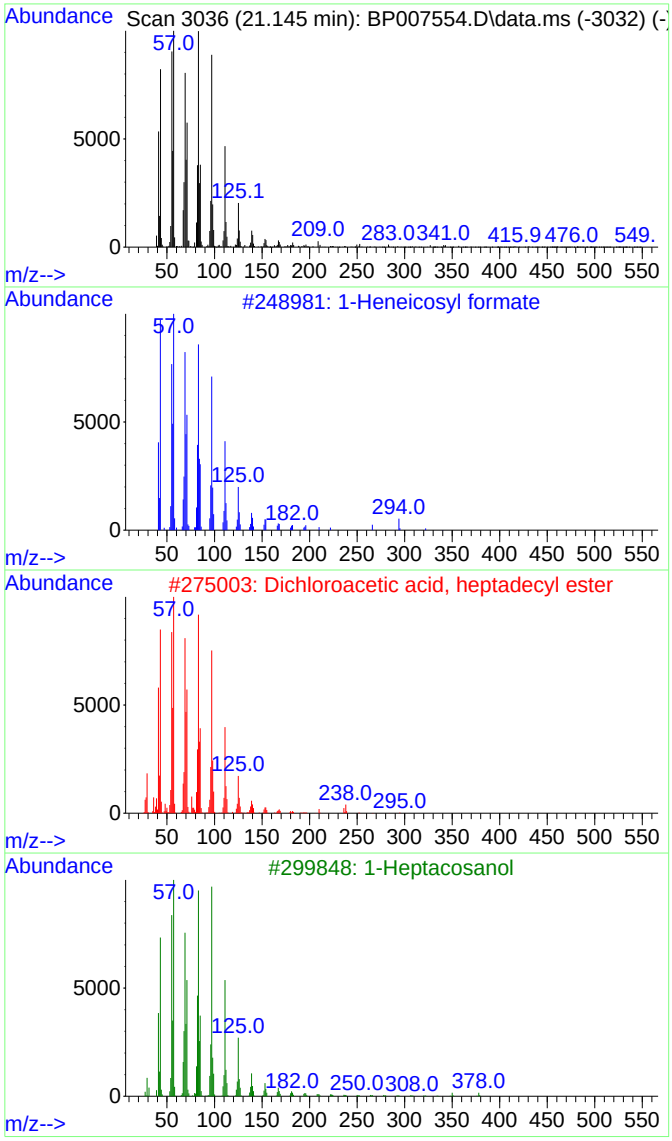
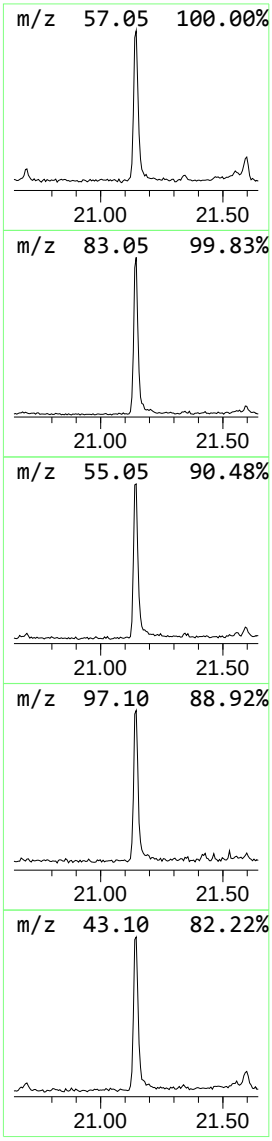
Peak Number 11 1-Heneicosyl formate Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.145	5.01 ng/ul	638321	Chrysene-d12	21.422

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Heneicosyl formate	340	C22H44O2	077899-03-7	97
2	Dichloroacetic acid, heptadecyl ...	366	C19H36Cl2O2	1000282-98-2	95
3	1-Heptacosanol	396	C27H56O	002004-39-9	94
4	Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	94
5	2- Chloropropionic acid, octadec...	360	C21H41ClO2	088104-31-8	94





Tentatively Identified Compound (LSC) summary

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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
Butanoic acid, ...	4.799	2.3	ng/ul	107580	1	7.952	943495	20.0
Butanoic acid, ...	4.952	3.0	ng/ul	140385	1	7.952	943495	20.0
Butanoic acid	5.393	2.0	ng/ul	95277	1	7.952	943495	20.0
Pentanoic acid,...	6.340	3.5	ng/ul	167216	1	7.952	943495	20.0
Hexanoic acid	6.952	2.3	ng/ul	108185	1	7.952	943495	20.0
Ethanol, 2-(2-e...	7.552	3.1	ng/ul	147477	1	7.952	943495	20.0
Butane, 1-(1-me...	7.740	5.2	ng/ul	246137	1	7.952	943495	20.0
Ethanol, 2-(2-b...	10.540	11.8	ng/ul	782509	2	10.757	1327710	20.0
Hexathiane	15.140	6.2	ng/ul	585379	3	14.587	1884390	20.0
n-Hexadecanoic ...	18.233	4.5	ng/ul	490262	4	17.339	2174350	20.0
1-Heneicosyl fo...	21.145	5.0	ng/ul	638321	5	21.422	2550000	20.0