LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102921\
Data File : BP007784.D
Acq On : 29 Oct 2021 14:44
Operator : CG/JU
Sample : M4262-20DL 2X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : BFY53DL

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-BP102521.M
Title : SVOA CALIBRATION

Signal : TIC: BP007784.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.405	16	20	29	rVB	178043	249687	0.65%	0.172%
2	3.681	54	67	71	rBV3	63794	184201	0.48%	0.127%
3	3.793	84	86	97	rBV	112363	202119	0.53%	0.139%
4	4.111	115	140	143	rBV	443366	2006804	5.26%	1.381%
5	4.834	248	263	275	rBV3	110340	450762	1.18%	0.310%
6	5.058	275	301	304	rBV3	408684	1742413	4.56%	1.199%
7	5.534	355	382	385	rBV	467117	2232033	5.85%	1.536%
8	6.340	506	519	527	rBV	49920	139651	0.37%	0.096%
9	6.464	533	540	551	rVB2	54890	129141	0.34%	0.089%
10	7.122	614	652	655	rBV5	1017196	6410284	16.79%	4.411%
11	7.258	670	675	682	rVB	222718	352248	0.92%	0.242%
12	7.458	702	709	720	rVB	230032	379822	0.99%	0.261%
13	7.928	779	789	797	rBV	932281	1575921	4.13%	1.084%
14	8.016	797	804	812	rVB4	76497	171408	0.45%	0.118%
15	8.181	824	832	842	rBV	157475	294880	0.77%	0.203%
16	8.458	872	879	887	rBV	68528	117715	0.31%	0.081%
17	8.628	887	908	914	rBV2	823934	3092151	8.10%	2.128%
18	8.699	914	920	934	rVB	3119064	5020844	13.15%	3.455%
19	9.022	968	975	979	rBV2	23768	39291	0.10%	0.027%
20	9.087	979	986	992	rBV	263240	436283	1.14%	0.300%
21	9.246	1006	1013	1021	rBV2	42062	82975	0.22%	0.057%
22	9.416	1038	1042	1048	rVV3	17591	41632	0.11%	0.029%
23	9.481	1048	1053	1058	rVV3	26444	50903	0.13%	0.035%
24	9.557	1058	1066	1070	rVV	169396	414612	1.09%	0.285%
25	9.599	1070	1073	1081	rVV2	101206	191975	0.50%	0.132%
26	9.810	1102	1109	1122	rVB	205460	355190	0.93%	0.244%
27	10.310	1143	1194	1197	rBV	2934865	23218238	60.82%	15.976%
28	10.352	1197	1201	1209	rVB	357221	641448	1.68%	0.441%
29	10.522	1223	1230	1239	rVB	294995	464508	1.22%	0.320%
30	10.728	1259	1265	1274	rVB	909263	1558920	4.08%	1.073%
31	10.869	1282	1289	1299	rBV	228796	421254	1.10%	0.290%
32	11.163	1333	1339	1344	rVB3	30246	49235	0.13%	0.034%
33	11.587	1400	1411	1415	rBV	577148	1157367	3.03%	0.796%
34	11.628	1415	1418	1428	rVV	72851	144117	0.38%	0.099%
35	12.222	1507	1519	1529	rBV	105963	205811	0.54%	0.142%
36	12.322	1529	1536	1542	rVV	426690	772335	2.02%	0.531%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102921\
Data File : BP007784.D
Acq On : 29 Oct 2021 14:44
Operator : CG/JU
Sample : M4262-20DL 2X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : BFY53DL

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-BP102521.M
Title : SVOA CALIBRATION

37	12.399	1542	1549	1553	rVV5	110518	300758	0.79%	0.207%
38	12.451	1553	1558	1570	rVV	206566	378701	0.99%	0.261%
39	12.546	1570	1574	1586	rVB7	29135	66320	0.17%	0.046%
40	12.828	1615	1622	1626	rBV	345537	684636	1.79%	0.471%
41	12.940	1626	1641	1659	rVB	2768852	9000111	23.58%	6.193%
42	13.175	1676	1681	1691	rVB	18211	42214	0.11%	0.029%
43	13.416	1717	1722	1729	rBV	109896	186358	0.49%	0.128%
44	13.722	1765	1774	1783	rBV	481885	815609	2.14%	0.561%
45	13.975	1811	1817	1831	rVV	705188	1039268	2.72%	0.715%
46	14.210	1852	1857	1859	rBV3	41067	64575	0.17%	0.044%
47	14.257	1859	1865	1875	rVB	771323	1138815	2.98%	0.784%
48	14.481	1888	1903	1906	rBV5	63177	166009	0.43%	0.114%
49	14.516	1906	1909	1912	rVV5	72355	129193	0.34%	0.089%
50	14.569	1912	1918	1929	rVV	2294322	3550583	9.30%	2.443%
51	14.663	1929	1934	1945	rVV6	143993	426703	1.12%	0.294%
52	14.763	1947	1951	1956	rVB	68100	112828	0.30%	0.078%
53	14.904	1969	1975	1977	rBV2	135833	224959	0.59%	0.155%
54	14.934	1977	1980	1985	rVV2	154394	249742	0.65%	0.172%
55	15.004	1986	1992	2002	rVB	931403	1355288	3.55%	0.933%
56	15.116	2006	2011	2015	rBV	73254	138275	0.36%	0.095%
57	15.210	2024	2027	2032	rVB	59867	79914	0.21%	0.055%
58	15.292	2036	2041	2052	rVB	218328	354193	0.93%	0.244%
59	15.422	2059	2063	2068	rVB	153966	209089	0.55%	0.144%
60	15.557	2080	2086	2098	rVB	1028684	1554456	4.07%	1.070%
61	15.675	2100	2106	2113	rBV	158098	239947	0.63%	0.165%
62	15.781	2121	2124	2131	rBV2	20361	41703	0.11%	0.029%
63	16.292	2208	2211	2218	rBV	76991	101001	0.26%	0.069%
64	16.351	2218	2221	2227	rVB	46155	58838	0.15%	0.040%
65	16.439	2230	2236	2245	rBV	24028	78200	0.20%	0.054%
66	16.516	2245	2249	2256	rVV5	35488	61410	0.16%	0.042%
67	16.745	2281	2288	2312	rVB	2932897	4442217	11.64%	3.057%
68	17.016	2331	2334	2341	rVB2	29945	41570	0.11%	0.029%
69	17.081	2341	2345	2351	rBV4	20048	42212	0.11%	0.029%
70	17.263	2371	2376	2380	rBV	180988	284365	0.74%	0.196%
71	17.322	2380	2386	2396	rVV	1947023	2830920	7.42%	1.948%
72	17.422	2397	2403	2409	rVB	1150045	1649353	4.32%	1.135%
73	17.504	2414	2417	2424	rVV5	40470	63469	0.17%	0.044%
74	17.722	2450	2454	2457	rBV	69297	100567	0.26%	0.069%
75	17.863	2474	2478	2481	rBV	231698	264091	0.69%	0.182%
76	17.904	2481	2485	2494	rVB	271618	357602	0.94%	0.246%

LSC Area Percent Report

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Data File : BP007784.D
Acq On : 29 Oct 2021 14:44
Operator : CG/JU
Sample : M4262-20DL 2X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : BFY53DL

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-BP102521.M
Title : SVOA CALIBRATION

77	18.004	2498	2502	2507	rBV2	61645	80424	0.21%	0.055%
78	18.269	2534	2547	2569	rBV	14090789	38175208	100.00%	26.267%
79	19.051	2677	2680	2683	rBV3	70440	78949	0.21%	0.054%
80	19.375	2730	2735	2741	rBV5	215081	490507	1.28%	0.338%
81	19.463	2745	2750	2771	rVB	4642373	6502987	17.03%	4.474%
82	19.657	2778	2783	2788	rVB	1469859	1933105	5.06%	1.330%
83	20.504	2924	2927	2939	rVB4	117636	225200	0.59%	0.155%
84	21.127	3030	3033	3042	rVB	204959	262005	0.69%	0.180%
85	21.327	3064	3067	3072	rBV	158638	169968	0.45%	0.117%
86	21.410	3076	3081	3087	rBV	2558149	3419843	8.96%	2.353%
87	23.127	3370	3373	3378	rVB	141472	196344	0.51%	0.135%
88	23.698	3463	3470	3477	rVB2	1058165	2140341	5.61%	1.473%
89	23.857	3490	3497	3510	rVB2	1758985	3735598	9.79%	2.570%

Sum of corrected areas: 145334719

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102921\
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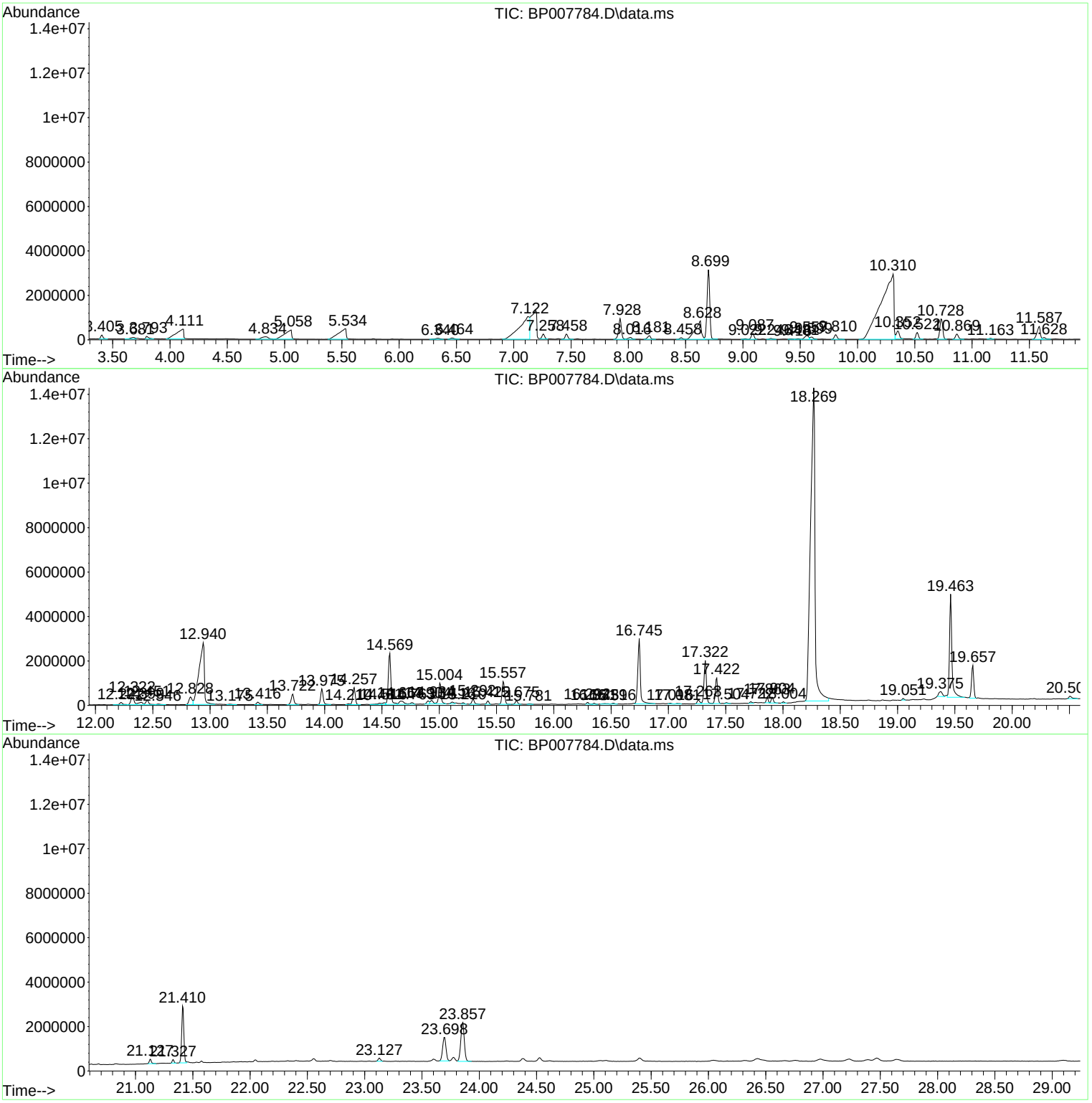
Acq On : 29 Oct 2021 14:44
Operator : CG/JU
Sample : M4262-20DL 2X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BFY53DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP102521.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102921\
Data File : BP007784.D
Acq On : 29 Oct 2021 14:44
Operator : CG/JU
Sample : M4262-20DL 2X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : BFY53DL

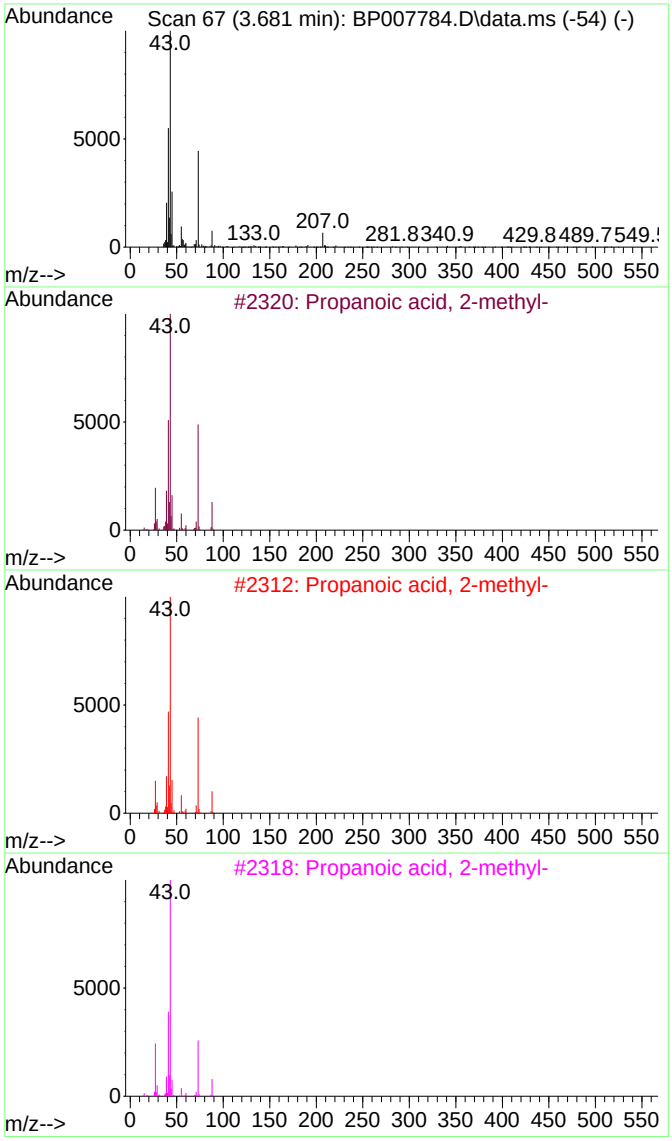
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP102521.M
Quant Title : SVOA CALIBRATION

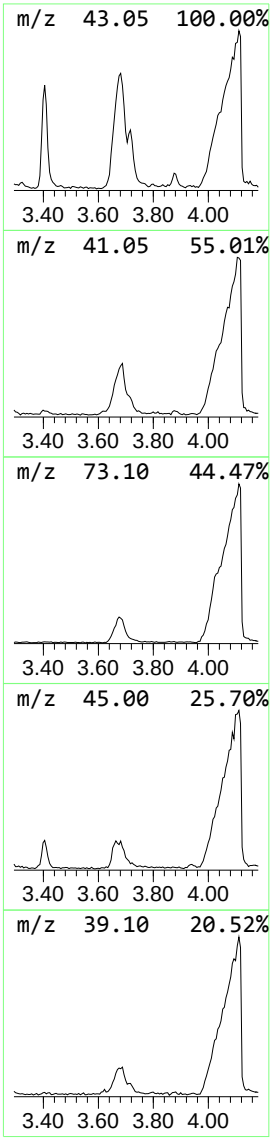
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 Propanoic acid, 2-methyl- Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.681	2.34 ng/ul	184201	1,4-Dichlorobenzene-d4	7.928

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Propanoic acid, 2-methyl-	88	C4H8O2	000079-31-2	52
2	Propanoic acid, 2-methyl-	88	C4H8O2	000079-31-2	50
3	Propanoic acid, 2-methyl-	88	C4H8O2	000079-31-2	47
4	Propanoic acid, 2-methyl-	88	C4H8O2	000079-31-2	47
5	Propanoic acid, 2-methyl-	88	C4H8O2	000079-31-2	30





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102921\
Data File : BP007784.D
Acq On : 29 Oct 2021 14:44
Operator : CG/JU
Sample : M4262-20DL 2X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BFY53DL

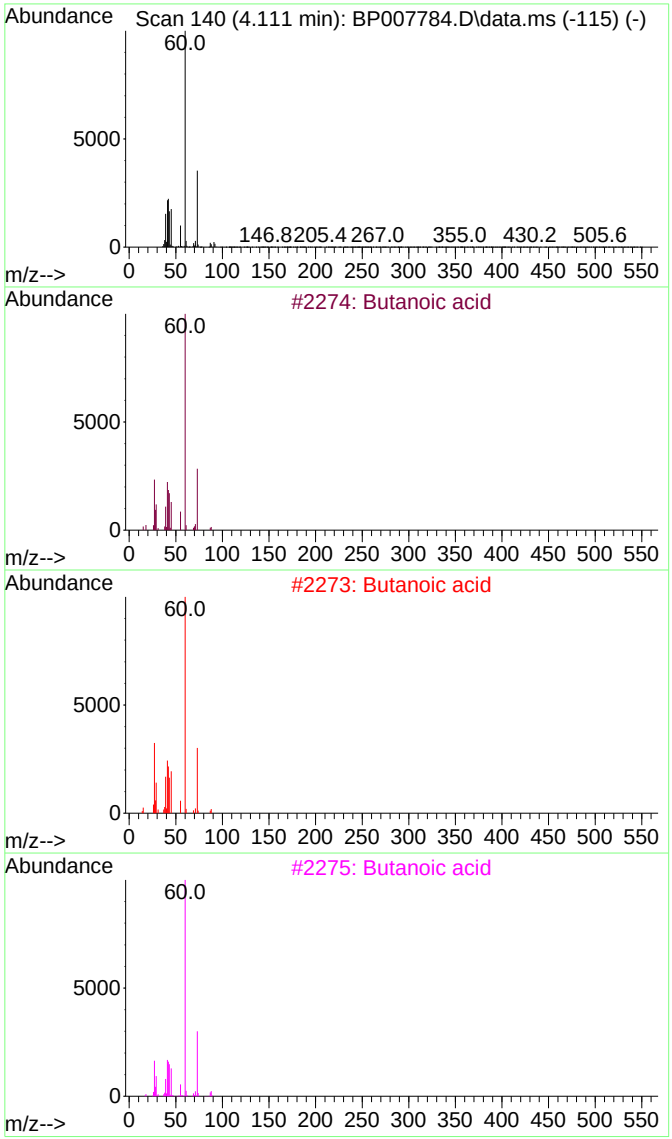
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP102521.M
Quant Title : SVOA CALIBRATION

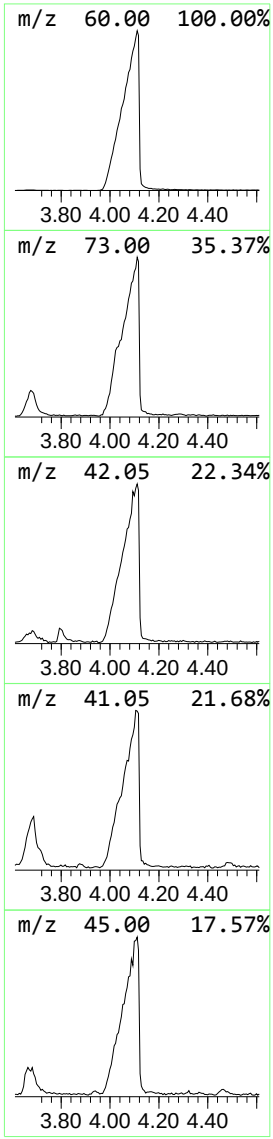
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 Butanoic acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.111	25.47 ng/ul	2006800	1,4-Dichlorobenzene-d4	7.928

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butanoic acid	88	C4H8O2	000107-92-6	91
2	Butanoic acid	88	C4H8O2	000107-92-6	86
3	Butanoic acid	88	C4H8O2	000107-92-6	72
4	Pentanoic acid	102	C5H10O2	000109-52-4	72
5	Butanoic acid	88	C4H8O2	000107-92-6	64





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102921\
Data File : BP007784.D
Acq On : 29 Oct 2021 14:44
Operator : CG/JU
Sample : M4262-20DL 2X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BFY53DL

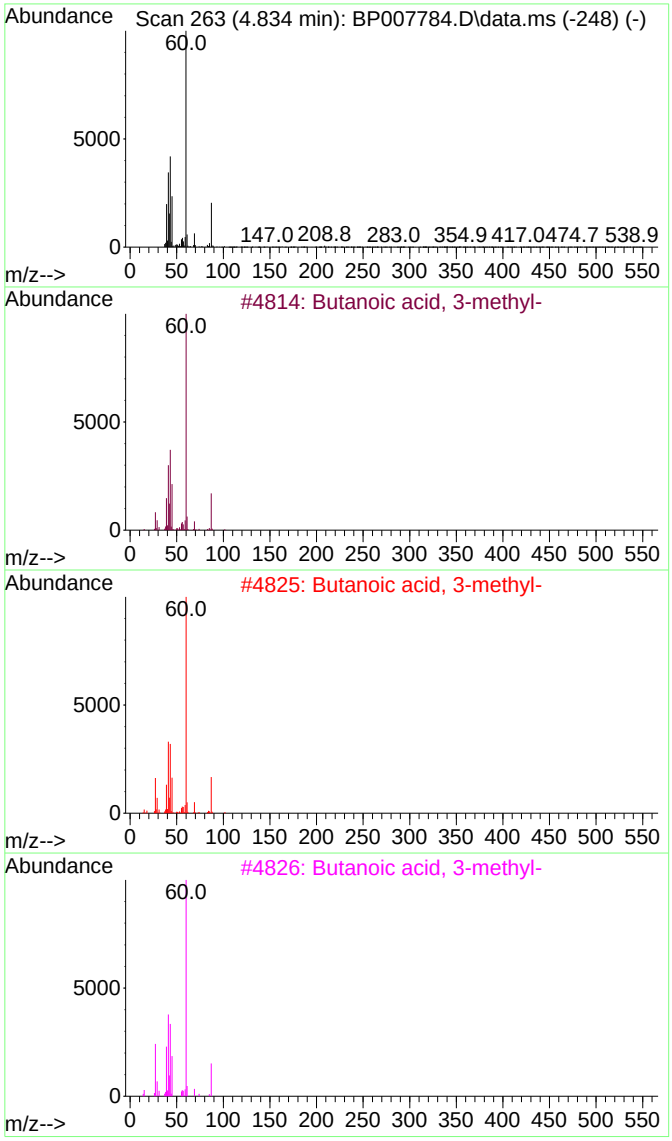
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP102521.M
Quant Title : SVOA CALIBRATION

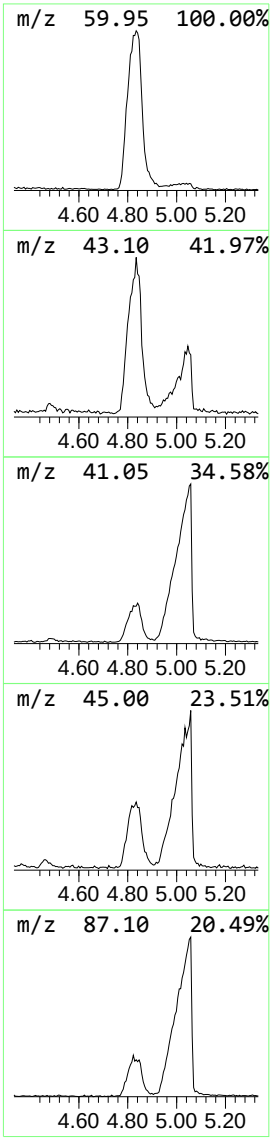
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Butanoic acid, 3-methyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.834	5.72 ng/ul	450762	1,4-Dichlorobenzene-d4	7.928

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butanoic acid, 3-methyl-	102	C5H10O2	000503-74-2	90
2	Butanoic acid, 3-methyl-	102	C5H10O2	000503-74-2	72
3	Butanoic acid, 3-methyl-	102	C5H10O2	000503-74-2	64
4	Butanoic acid, 3-methyl-	102	C5H10O2	000503-74-2	64
5	Butanoic acid, 3-methyl-	102	C5H10O2	000503-74-2	64





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102921\
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Acq On : 29 Oct 2021 14:44
Operator : CG/JU
Sample : M4262-20DL 2X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : BFY53DL

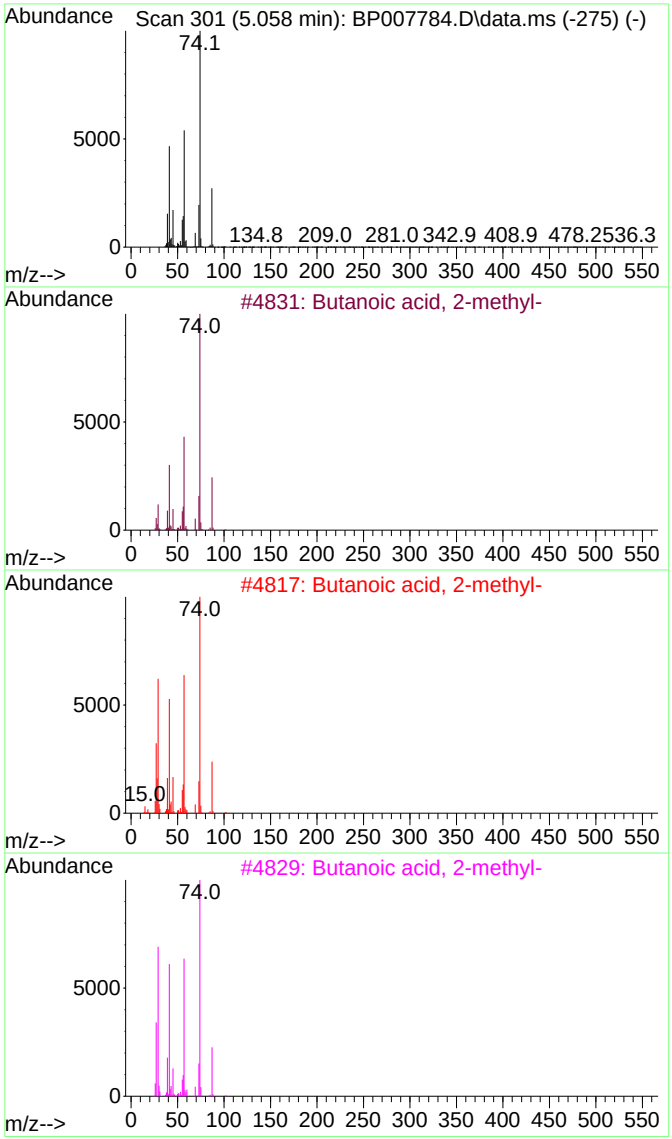
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP102521.M
Quant Title : SVOA CALIBRATION

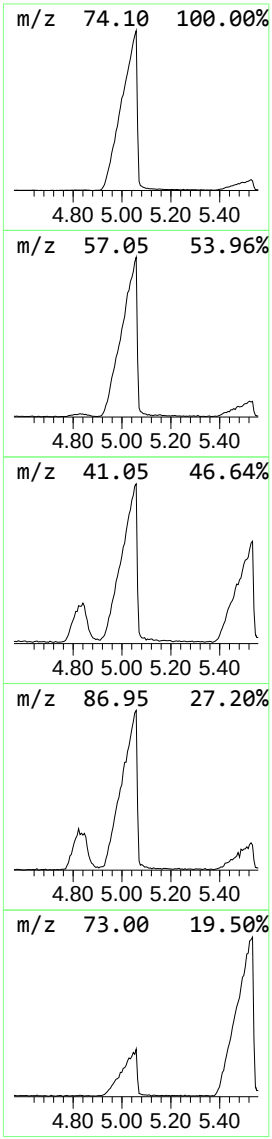
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Butanoic acid, 2-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.058	22.11 ng/ul	1742410	1,4-Dichlorobenzene-d4	7.928

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	78
3	Butanoic acid, 2-methyl-	102	C5H10O2	000116-53-0	78
4	Hexanoic acid, 2-methyl-	130	C7H14O2	004536-23-6	64
5	Hexanoic acid, 2-methyl-	130	C7H14O2	004536-23-6	64





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102921\
Data File : BP007784.D
Acq On : 29 Oct 2021 14:44
Operator : CG/JU
Sample : M4262-20DL 2X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BFY53DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP102521.M
Quant Title : SVOA CALIBRATION

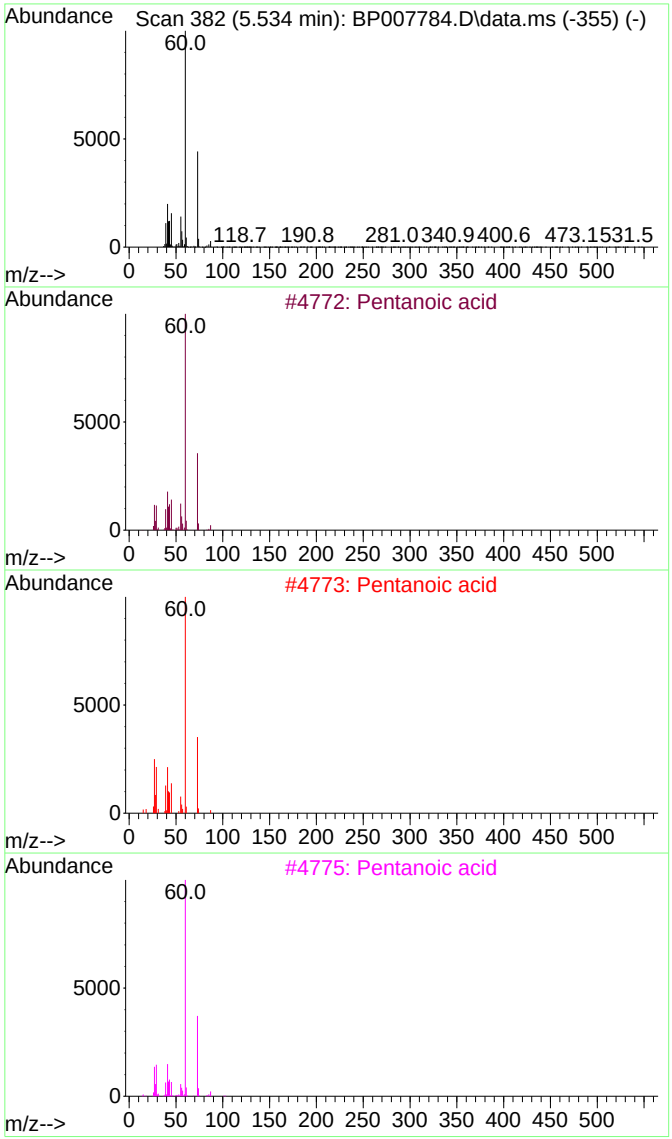
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

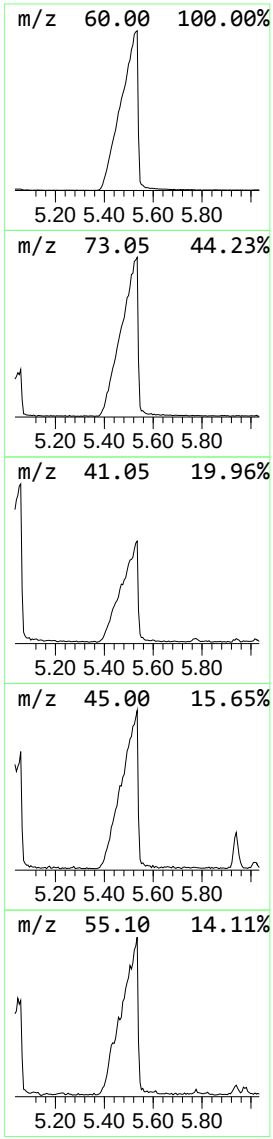
****

Peak Number 5 Pentanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.534	28.33 ng/ul	2232030	1,4-Dichlorobenzene-d4	7.928

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentanoic acid	102	C5H10O2	000109-52-4	90
2	Pentanoic acid	102	C5H10O2	000109-52-4	83
3	Pentanoic acid	102	C5H10O2	000109-52-4	78
4	Propanedioic acid, propyl-	146	C6H10O4	000616-62-6	74
5	Butanoic acid	88	C4H8O2	000107-92-6	72





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102921\
Data File : BP007784.D
Acq On : 29 Oct 2021 14:44
Operator : CG/JU
Sample : M4262-20DL 2X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : BFY53DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP102521.M
Quant Title : SVOA CALIBRATION

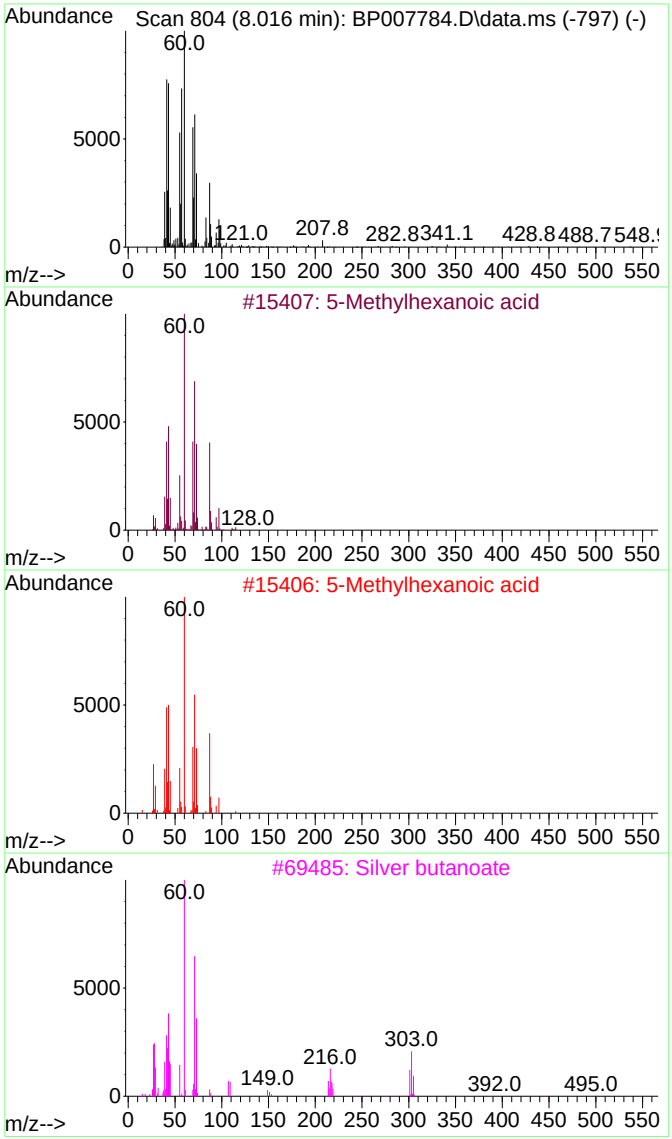
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

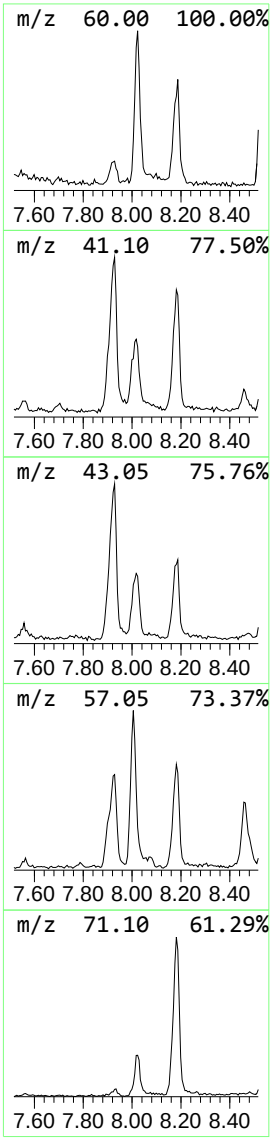
****

Peak Number 6 5-Methylhexanoic acid Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.016	2.18 ng/ul	171408	1,4-Dichlorobenzene-d4	7.928

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	5-Methylhexanoic acid	130	C7H14O2	000628-46-6	58
2	5-Methylhexanoic acid	130	C7H14O2	000628-46-6	53
3	Silver butanoate	194	C4H7AgO2	005076-24-4	40
4	.beta.-D-Glucopyranose, 1,6-anhy...	162	C6H10O5	000498-07-7	40
5	Butanoic acid	88	C4H8O2	000107-92-6	38





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Acq On : 29 Oct 2021 14:44
Operator : CG/JU
Sample : M4262-20DL 2X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : BFY53DL

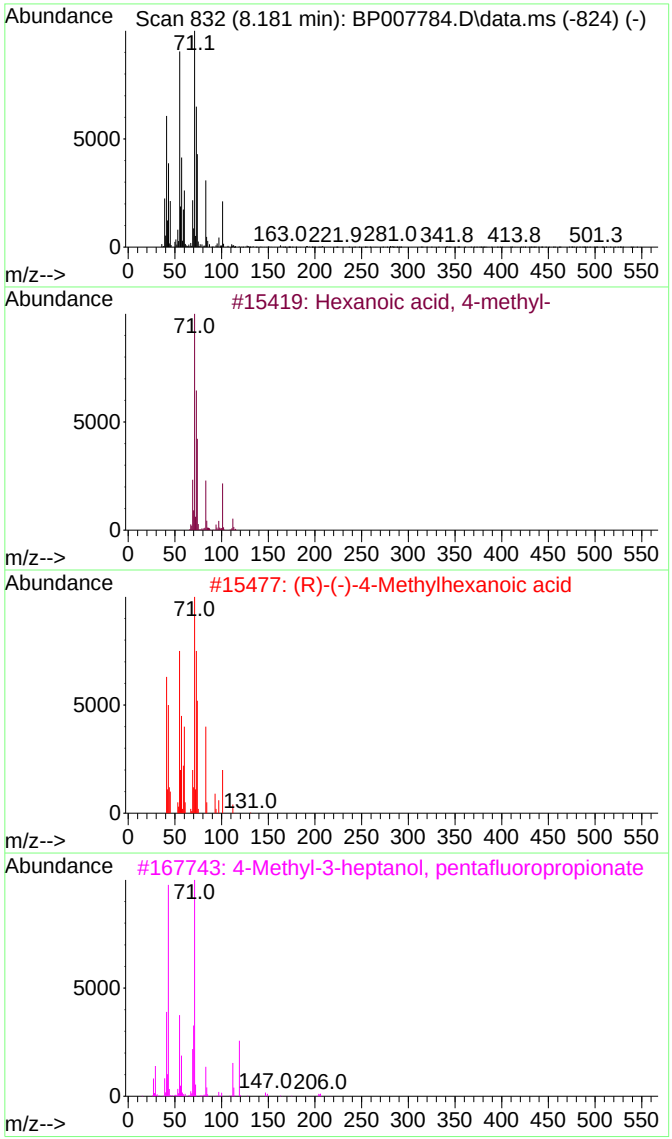
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP102521.M
Quant Title : SVOA CALIBRATION

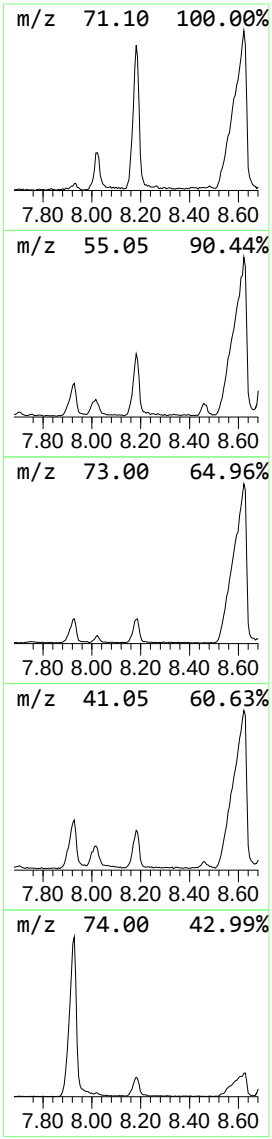
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 Hexanoic acid, 4-methyl- Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.181	3.74 ng/ul	294880	1,4-Dichlorobenzene-d4	7.928

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexanoic acid, 4-methyl-	130	C7H14O2	001561-11-1	90
2	(R)-(-)-4-Methylhexanoic acid	130	C7H14O2	052745-93-4	50
3	4-Methyl-3-heptanol, pentafluoro...	276	C11H17F5O2	1000365-51-7	27
4	Butane, 1-bromo-2-methyl-, (S)-	150	C5H11Br	000534-00-9	22
5	Sulfurous acid, 2-pentyl undecyl...	306	C16H34O3S	1000309-16-0	16





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Acq On : 29 Oct 2021 14:44
Operator : CG/JU
Sample : M4262-20DL 2X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BFY53DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP102521.M
Quant Title : SVOA CALIBRATION

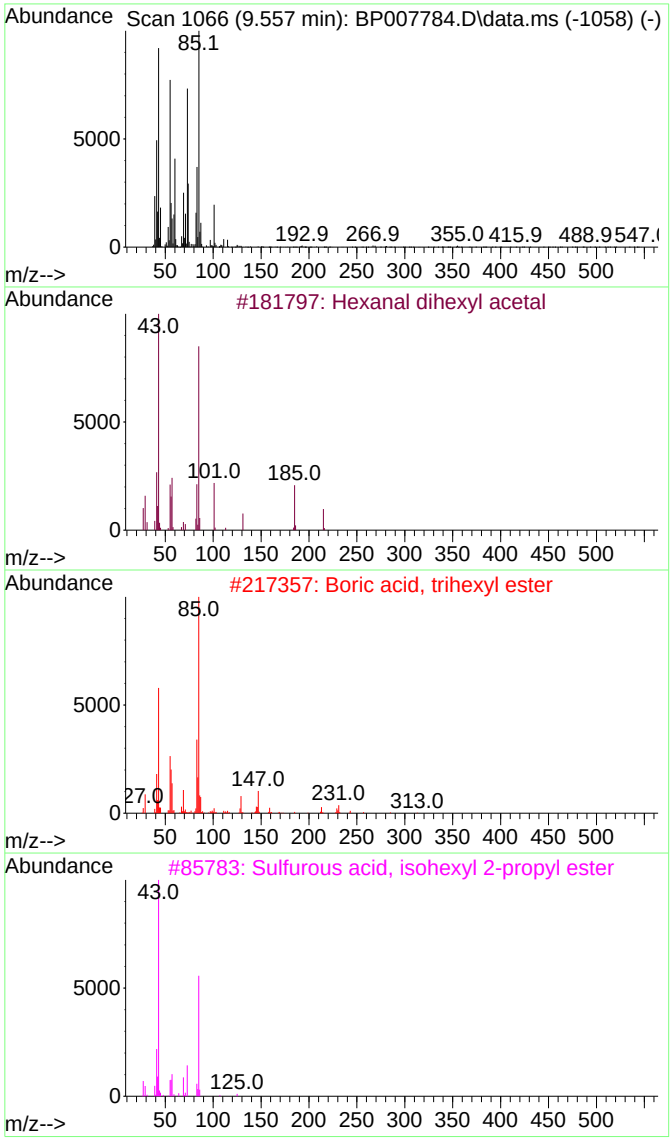
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

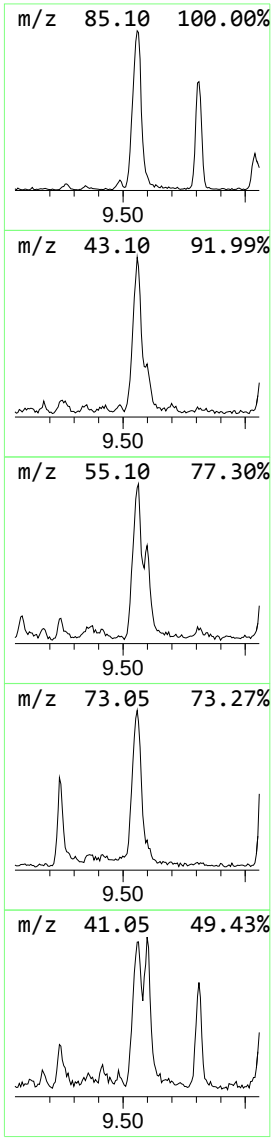
****

Peak Number 8 unknown-01 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.557	5.32 ng/ul	414612	Naphthalene-d8	10.728

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexanal dihexyl acetal	286	C18H38O2	1000430-99-9	43
2	Boric acid, trihexyl ester	314	C18H39BO3	005337-36-0	38
3	Sulfurous acid, isohexyl 2-propy...	208	C9H20O3S	1000309-11-6	38
4	Oxazole, 4,5-dihydro-2-methyl-	85	C4H7NO	001120-64-5	30
5	2-Butene, 1,1-dimethoxy-	116	C6H12O2	021962-24-3	30





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Acq On : 29 Oct 2021 14:44
Operator : CG/JU
Sample : M4262-20DL 2X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : BFY53DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP102521.M
Quant Title : SVOA CALIBRATION

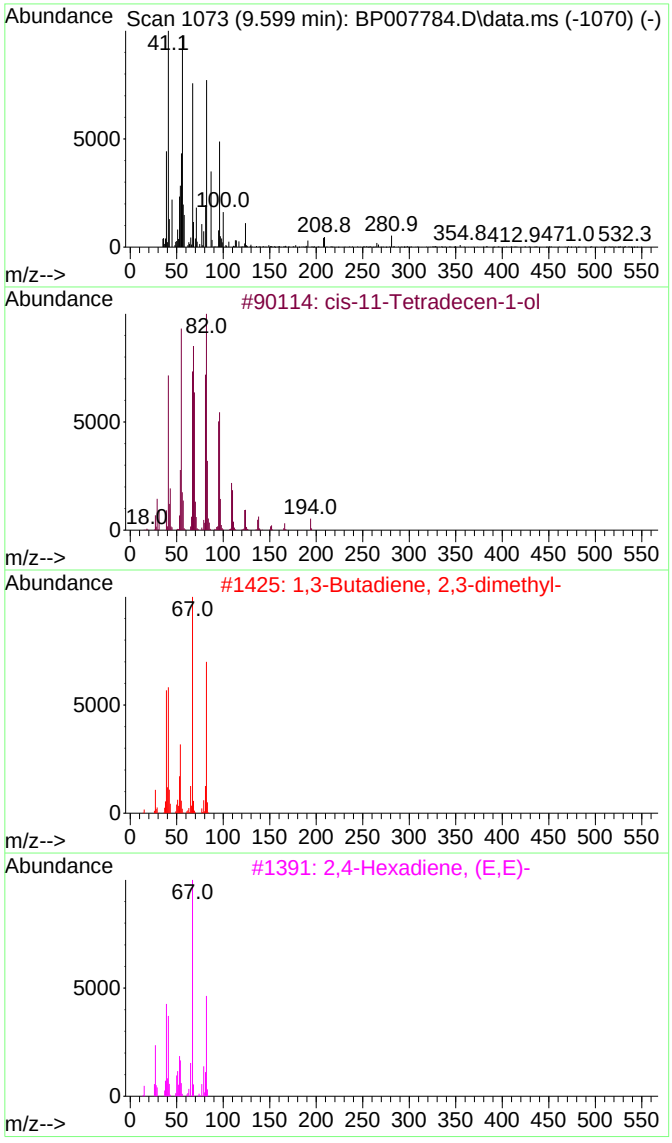
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

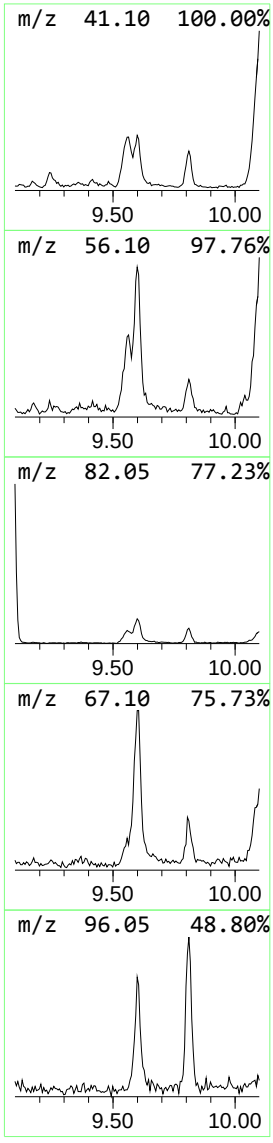
****

Peak Number 9 unknown-02 Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.599	2.46 ng/ul	191975	Naphthalene-d8	10.728

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	cis-11-Tetradecen-1-ol	212	C14H28O	034010-15-6	47
2	1,3-Butadiene, 2,3-dimethyl-	82	C6H10	000513-81-5	43
3	2,4-Hexadiene, (E,E)-	82	C6H10	005194-51-4	38
4	3-Hexyne	82	C6H10	000928-49-4	38
5	1,3-Pentadiene, 2-methyl-	82	C6H10	001118-58-7	38





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Data File : BP007784.D
Acq On : 29 Oct 2021 14:44
Operator : CG/JU
Sample : M4262-20DL 2X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : BFY53DL

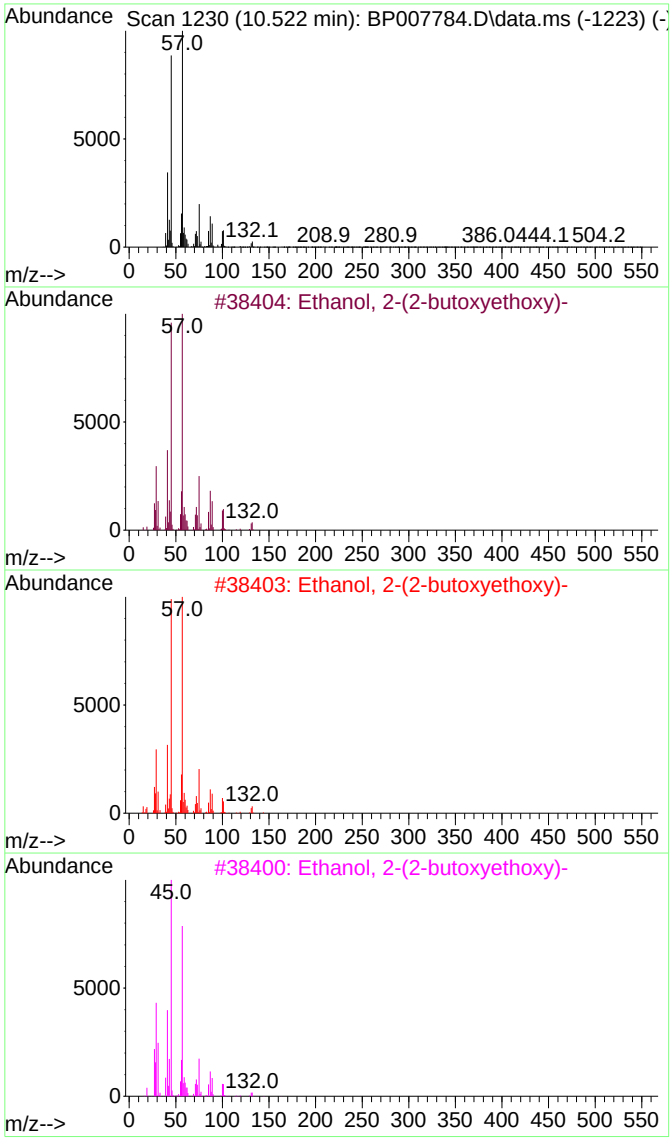
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP102521.M
Quant Title : SVOA CALIBRATION

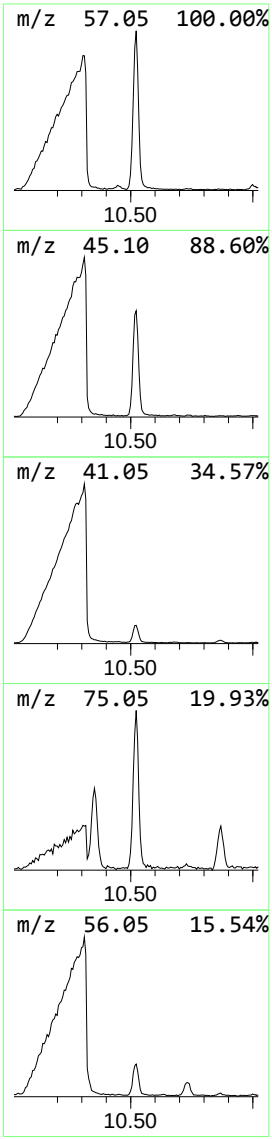
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 Ethanol, 2-(2-butoxyethoxy)- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.522	5.96 ng/ul	464508	Naphthalene-d8	10.728

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	91
2	Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
3	Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
4	Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
5	Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	72





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Acq On : 29 Oct 2021 14:44
Operator : CG/JU
Sample : M4262-20DL 2X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BFY53DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP102521.M
Quant Title : SVOA CALIBRATION

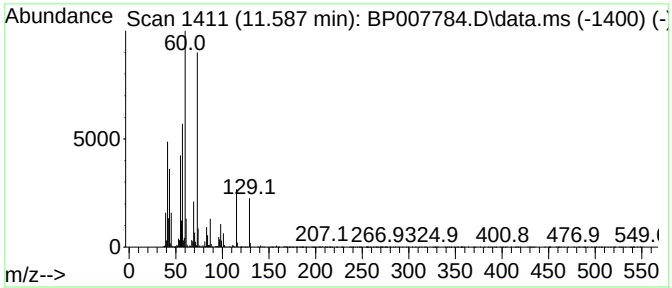
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

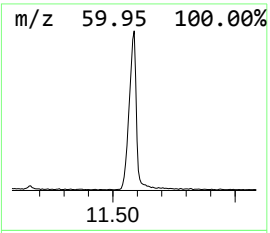
****

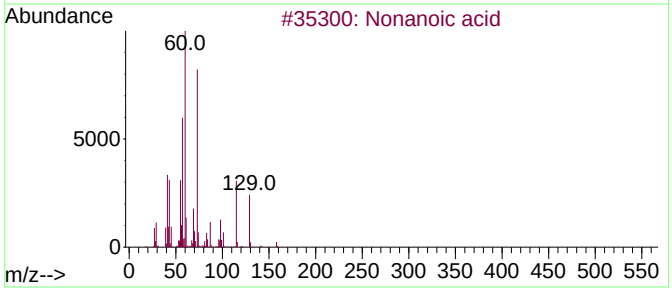
Peak Number 11 Nonanoic acid Concentration Rank 8

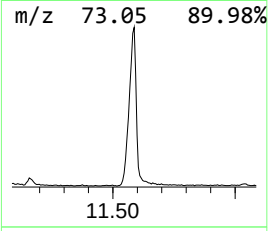
R.T.	EstConc	Area	Relative to ISTD	R.T.
11.587	14.85 ng/ul	1157370	Naphthalene-d8	10.728

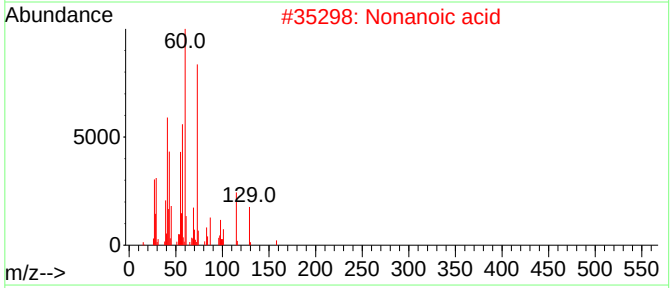
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonanoic acid	158	C9H18O2	000112-05-0	95
2	Nonanoic acid	158	C9H18O2	000112-05-0	86
3	Nonanoic acid	158	C9H18O2	000112-05-0	86
4	n-Decanoic acid	172	C10H20O2	000334-48-5	72
5	n-Decanoic acid	172	C10H20O2	000334-48-5	72

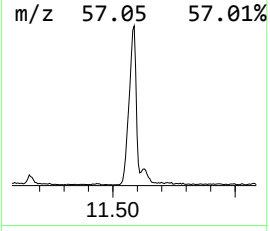


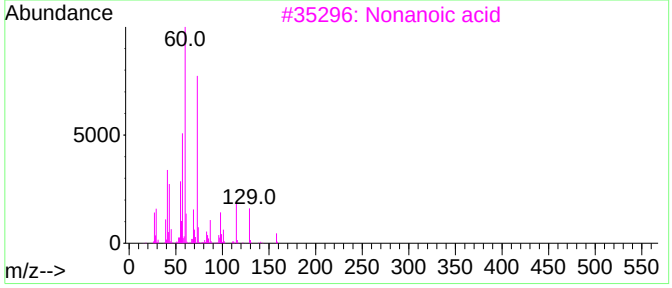


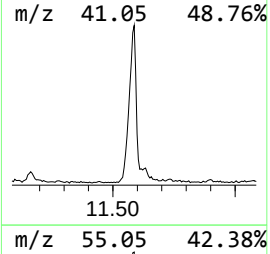


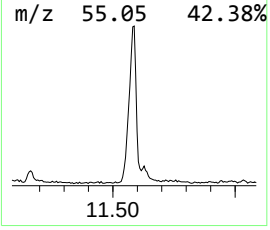












Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102921\
Data File : BP007784.D
Acq On : 29 Oct 2021 14:44
Operator : CG/JU
Sample : M4262-20DL 2X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : BFY53DL

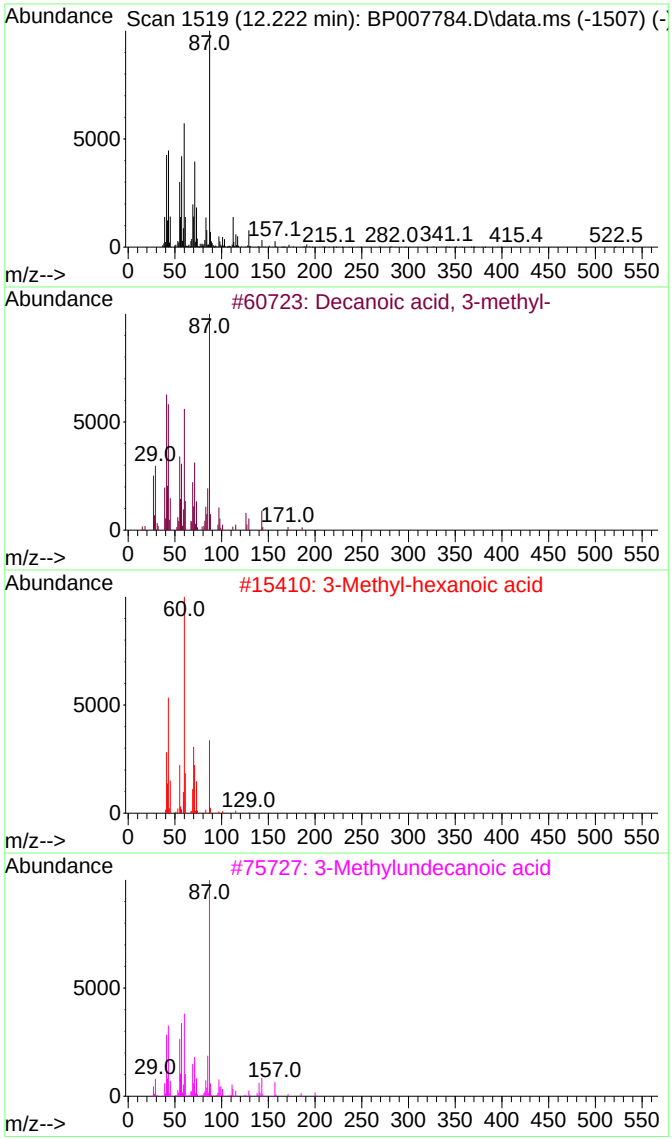
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP102521.M
Quant Title : SVOA CALIBRATION

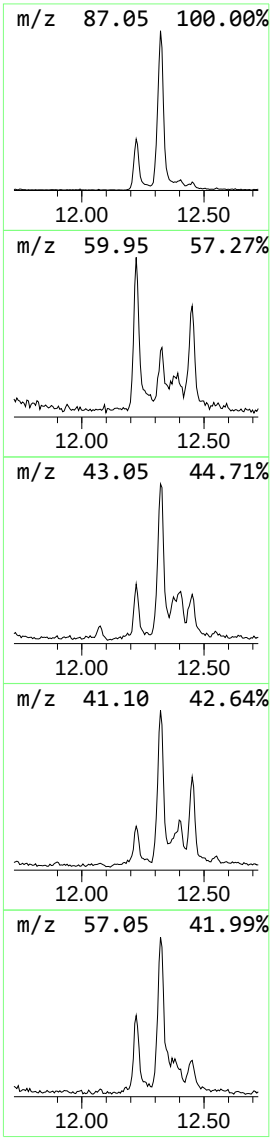
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 12 Decanoic acid, 3-methyl- Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.222	2.64 ng/ul	205811	Naphthalene-d8	10.728

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Decanoic acid, 3-methyl-	186	C11H22O2	060308-82-9	78
2	3-Methyl-hexanoic acid	130	C7H14O2	1000365-65-4	59
3	3-Methylundecanoic acid	200	C12H24O2	065781-38-6	53
4	3-Methyloctanoic acid	158	C9H18O2	006061-10-5	50
5	Pentanoic acid, 3-methyl-	116	C6H12O2	000105-43-1	40





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Acq On : 29 Oct 2021 14:44
Operator : CG/JU
Sample : M4262-20DL 2X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : BFY53DL

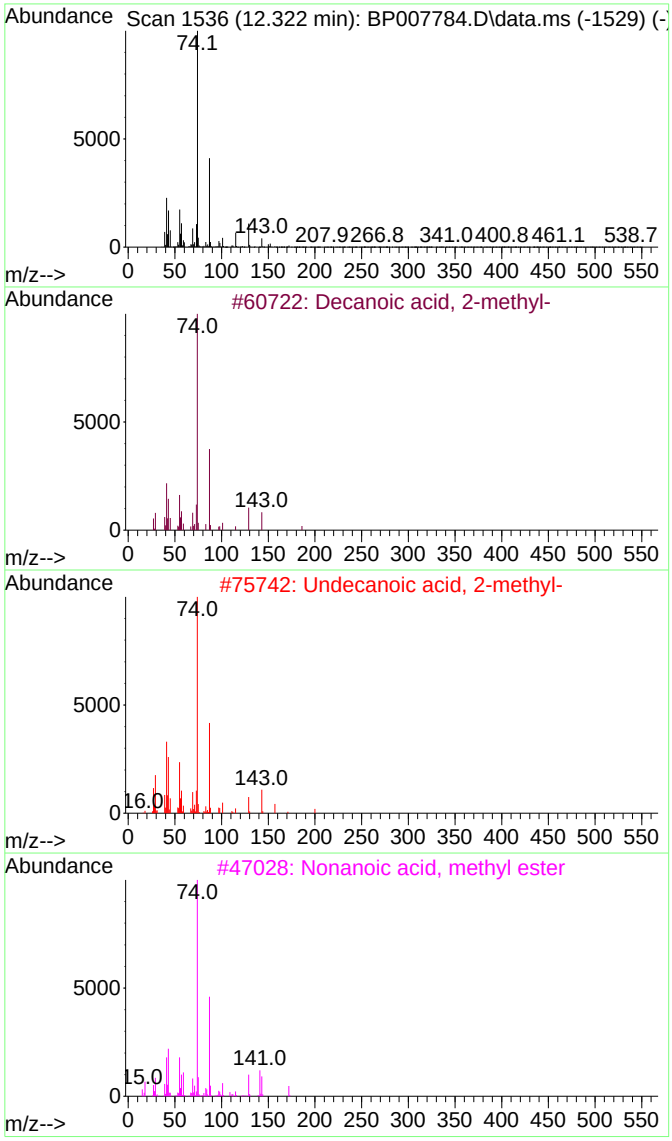
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP102521.M
Quant Title : SVOA CALIBRATION

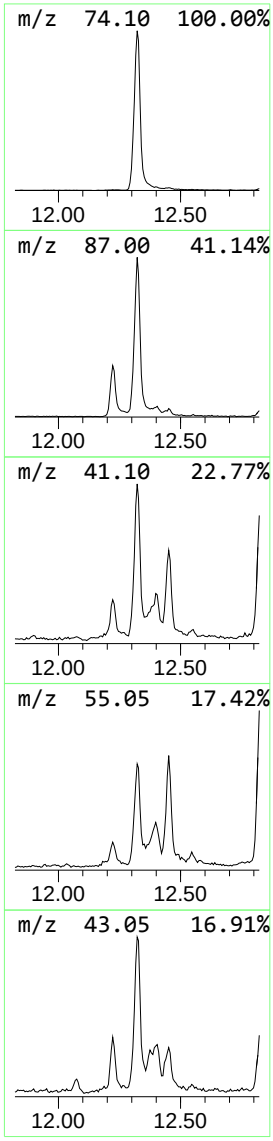
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 13 Decanoic acid, 2-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.322	9.91 ng/ul	772335	Naphthalene-d8	10.728

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Decanoic acid, 2-methyl-	186	C11H22O2	024323-23-7	86
2	Undecanoic acid, 2-methyl-	200	C12H24O2	024323-25-9	86
3	Nonanoic acid, methyl ester	172	C10H20O2	001731-84-6	72
4	Nonanoic acid, methyl ester	172	C10H20O2	001731-84-6	64
5	Nonanoic acid, methyl ester	172	C10H20O2	001731-84-6	56





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Acq On : 29 Oct 2021 14:44
Operator : CG/JU
Sample : M4262-20DL 2X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : BFY53DL

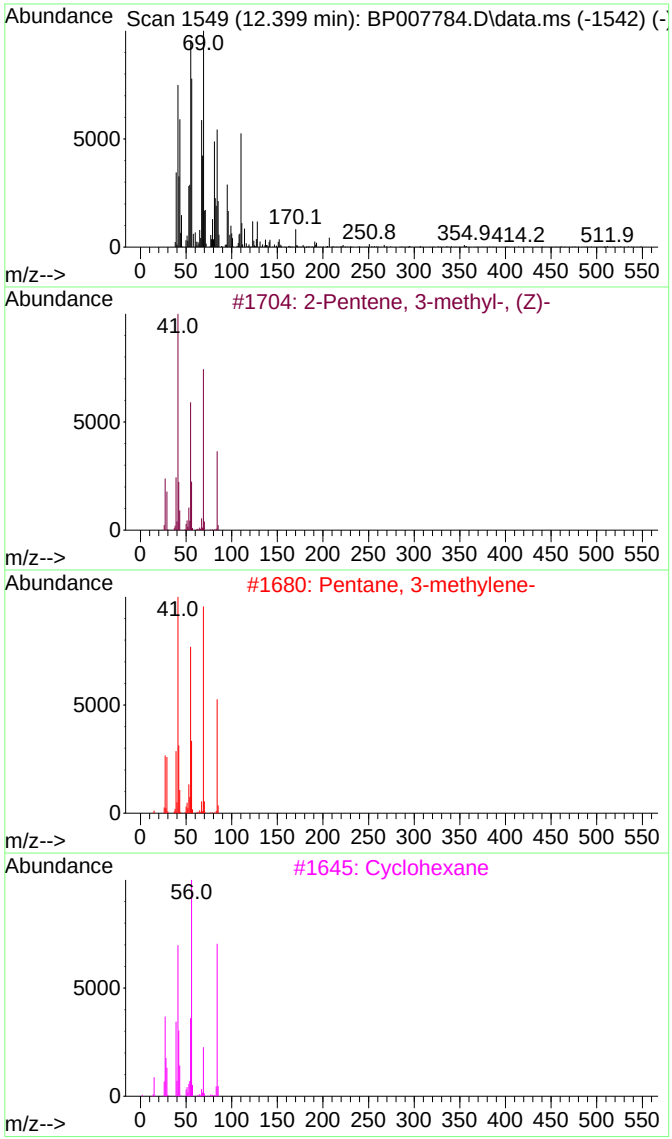
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP102521.M
Quant Title : SVOA CALIBRATION

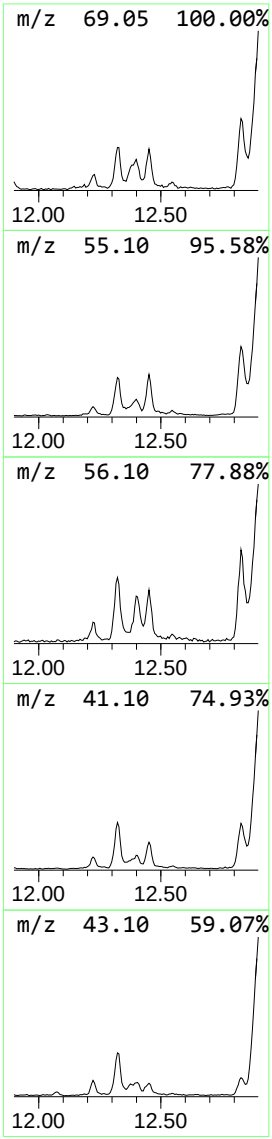
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 14 (DEL) Alkane: Straight-Chai... Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.399	3.86 ng/ul	300758	Naphthalene-d8	10.728

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentene, 3-methyl-, (Z)-	84	C6H12	000922-62-3	30
2	Pentane, 3-methylene-	84	C6H12	000760-21-4	27
3	Cyclohexane	84	C6H12	000110-82-7	22
4	3-Penten-2-one	84	C5H8O	000625-33-2	22
5	1,5-Heptadiene, 2,3,6-trimethyl-	138	C10H18	033501-88-1	14





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Acq On : 29 Oct 2021 14:44
Operator : CG/JU
Sample : M4262-20DL 2X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BFY53DL

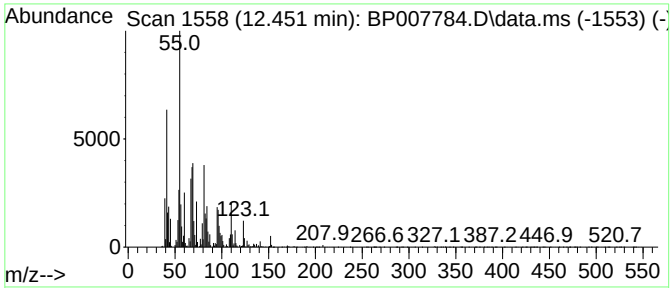
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP102521.M
Quant Title : SVOA CALIBRATION

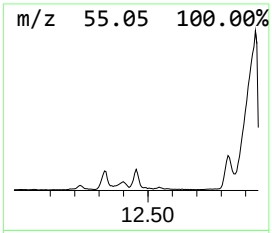
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

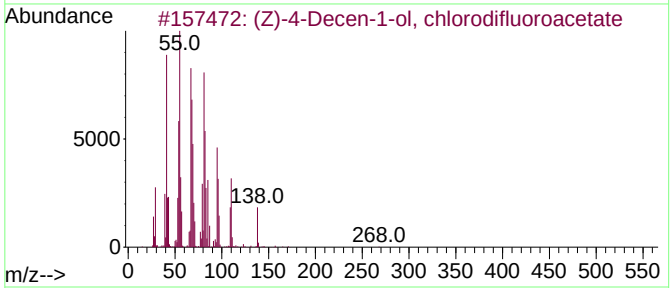
****
Peak Number 15 unknown-03 Concentration Rank 14

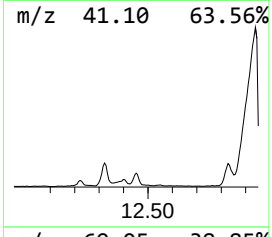
R.T.	EstConc	Area	Relative to ISTD	R.T.
12.451	4.86 ng/ul	378701	Naphthalene-d8	10.728

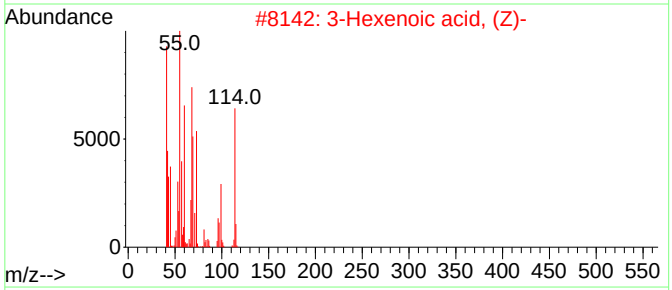
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	(Z)-4-Decen-1-ol, chlorodifluoro...	268	C12H19ClF2O2	1000447-45-6	46
2	3-Hexenoic acid, (Z)-	114	C6H10O2	001775-43-5	43
3	Z-12-Tetradecen-1-ol	212	C14H28O	1000130-84-0	38
4	3-Hexenoic acid, (E)-	114	C6H10O2	001577-18-0	35
5	11-Tetradecen-1-ol, (E)-	212	C14H28O	035153-18-5	27

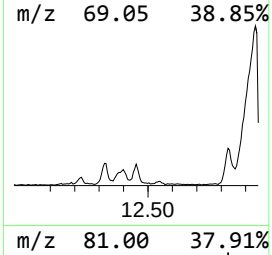


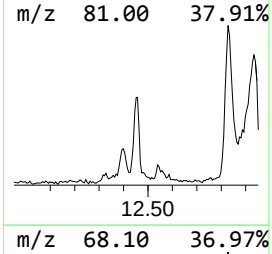


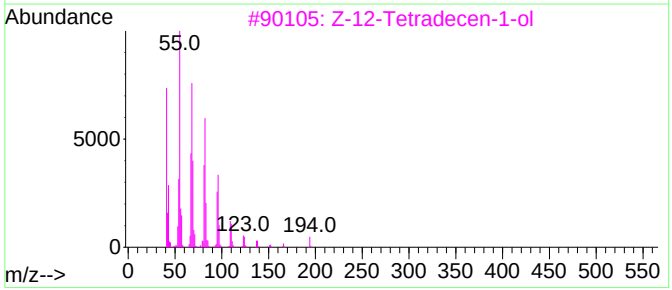


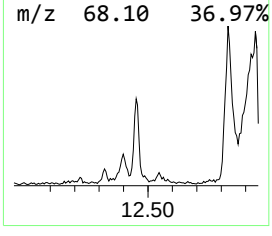












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Data File : BP007784.D
Acq On : 29 Oct 2021 14:44
Operator : CG/JU
Sample : M4262-20DL 2X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : BFY53DL

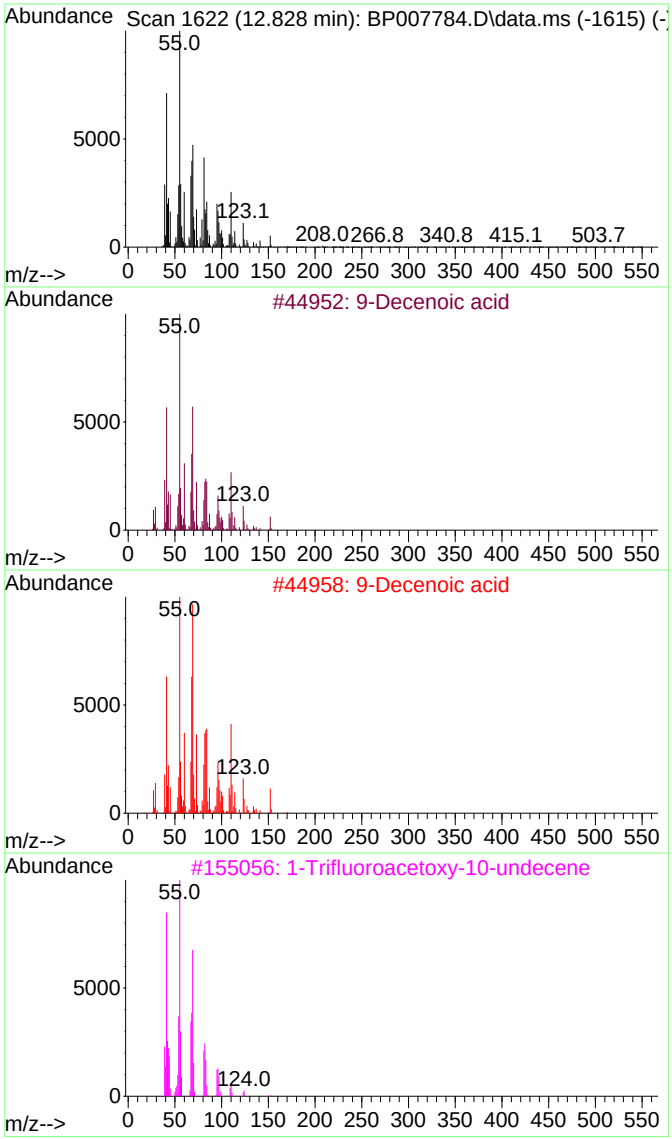
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP102521.M
Quant Title : SVOA CALIBRATION

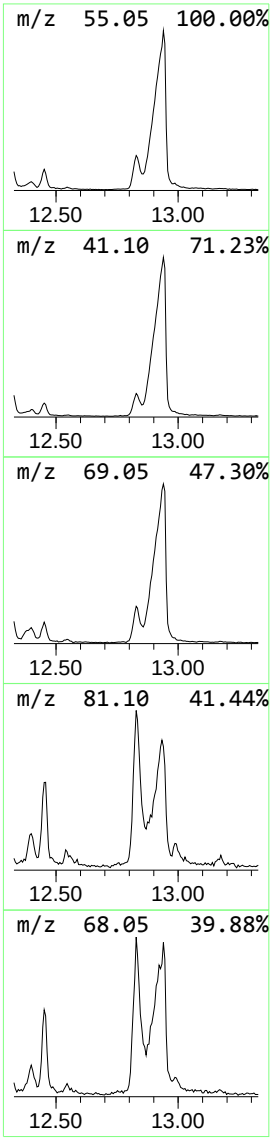
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 16 9-Decenoic acid Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.828	3.86 ng/ul	684636	Acenaphthene-d10	14.569

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9-Decenoic acid	170	C10H18O2	014436-32-9	64
2	9-Decenoic acid	170	C10H18O2	014436-32-9	43
3	1-Trifluoroacetoxy-10-undecene	266	C13H21F3O2	001675-20-3	41
4	3-Nonen-1-ol, (E)-	142	C9H18O	010339-61-4	38
5	11-Tetradecen-1-ol, (E)-	212	C14H28O	035153-18-5	38





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Operator : CG/JU
Sample : M4262-20DL 2X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : BFY53DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP102521.M
Quant Title : SVOA CALIBRATION

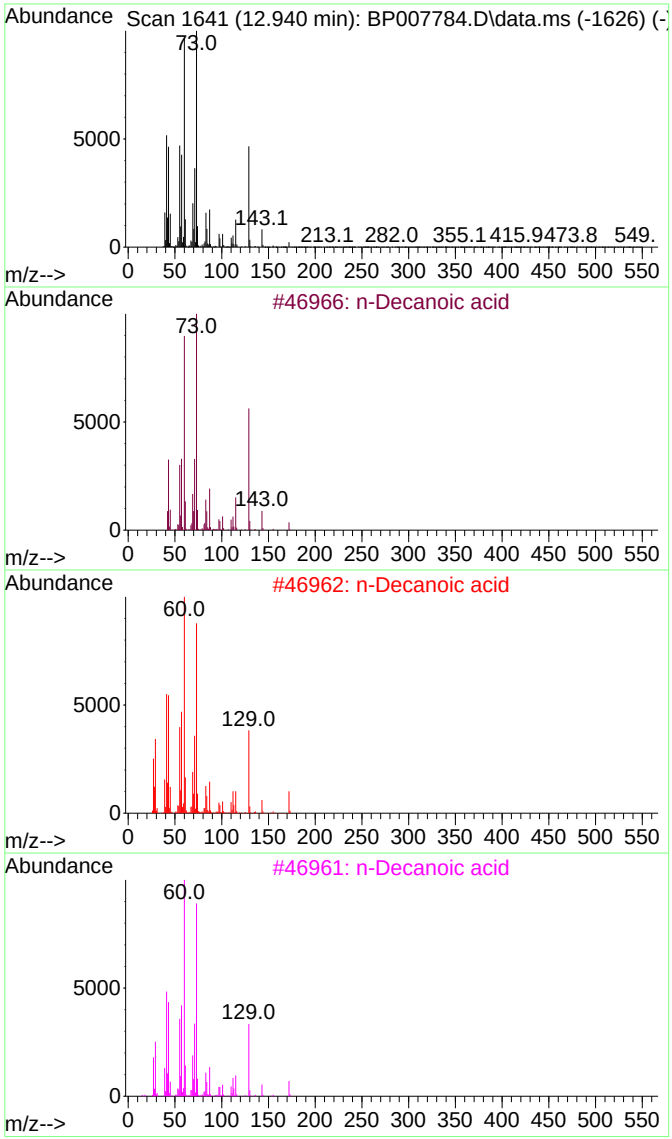
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

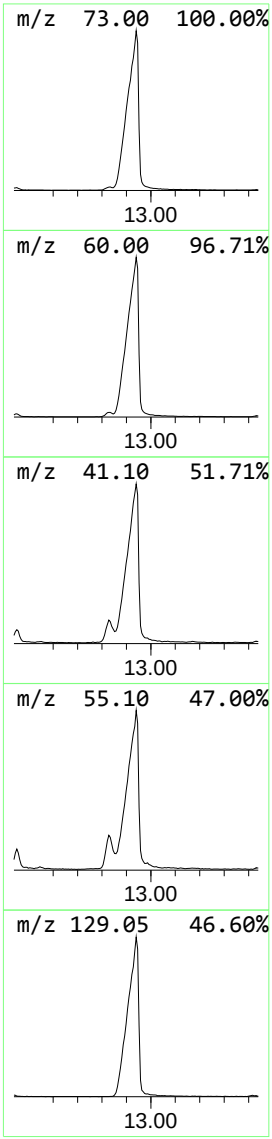
****

Peak Number 17 n-Decanoic acid Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.940	50.70 ng/ul	9000110	Acenaphthene-d10	14.569

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Decanoic acid	172	C10H20O2	000334-48-5	97
2	n-Decanoic acid	172	C10H20O2	000334-48-5	91
3	n-Decanoic acid	172	C10H20O2	000334-48-5	83
4	n-Decanoic acid	172	C10H20O2	000334-48-5	64
5	n-Decanoic acid	172	C10H20O2	000334-48-5	53





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Acq On : 29 Oct 2021 14:44
Operator : CG/JU
Sample : M4262-20DL 2X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : BFY53DL

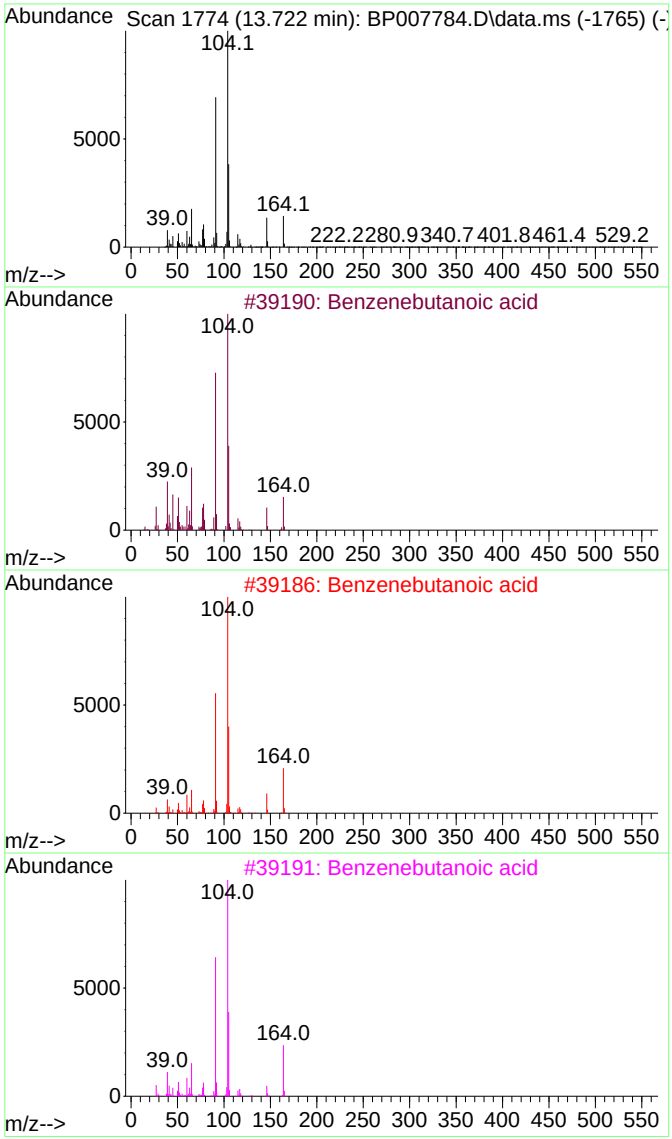
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP102521.M
Quant Title : SVOA CALIBRATION

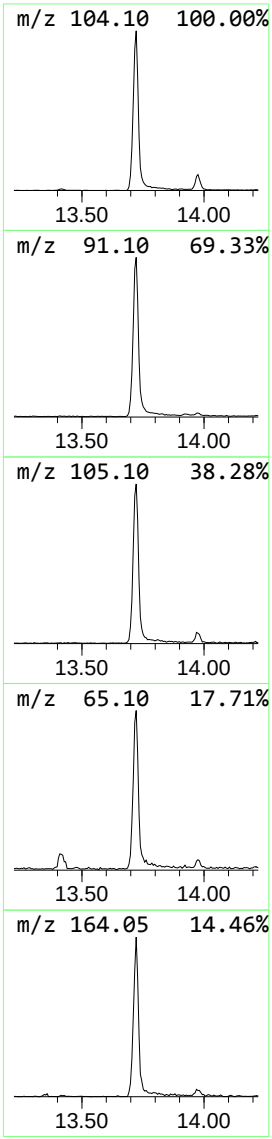
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 18 Benzenebutanoic acid Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.722	4.59 ng/ul	815609	Acenaphthene-d10	14.569

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzenebutanoic acid	164	C10H12O2	001821-12-1	91
2	Benzenebutanoic acid	164	C10H12O2	001821-12-1	90
3	Benzenebutanoic acid	164	C10H12O2	001821-12-1	87
4	2-Phenylethylamine, N,N-di(trifl...	313	C12H9F6NO2	1000463-67-7	72
5	Octanoic acid, 2-phenylethyl ester	248	C16H24O2	005457-70-5	64





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Data File : BP007784.D
Acq On : 29 Oct 2021 14:44
Operator : CG/JU
Sample : M4262-20DL 2X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : BFY53DL

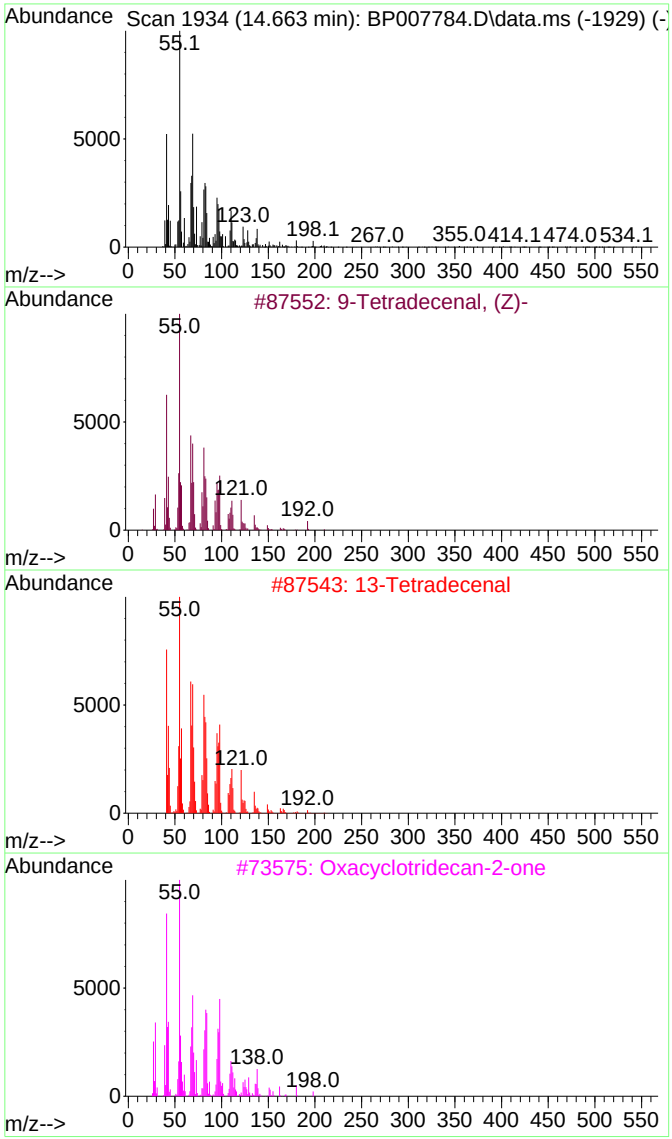
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP102521.M
Quant Title : SVOA CALIBRATION

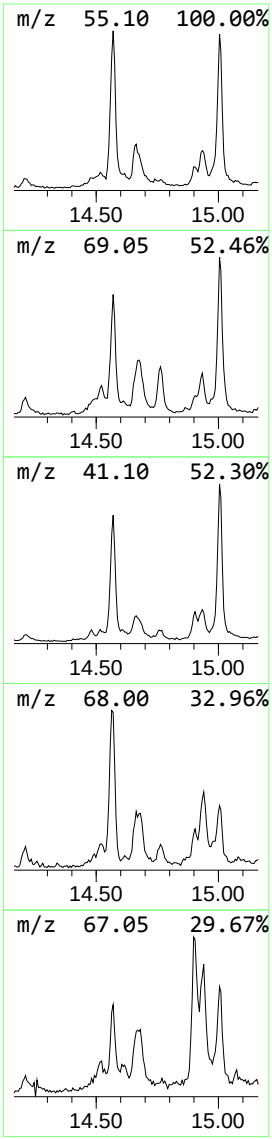
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 19 9-Tetradecenal, (Z)- Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.663	2.40 ng/ul	426703	Acenaphthene-d10	14.569

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9-Tetradecenal, (Z)-	210	C14H26O	053939-27-8	91
2	13-Tetradecenal	210	C14H26O	085896-31-7	87
3	Oxacyclotridecan-2-one	198	C12H22O2	000947-05-7	76
4	Undec-10-ynoic acid, dodecyl ester	350	C23H42O2	1010406-16-5	74
5	(E)-tetradec-11-en-1-yl 2,2,3,3,...	358	C17H27F5O2	1000385-85-0	68





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Acq On : 29 Oct 2021 14:44
Operator : CG/JU
Sample : M4262-20DL 2X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : BFY53DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP102521.M
Quant Title : SVOA CALIBRATION

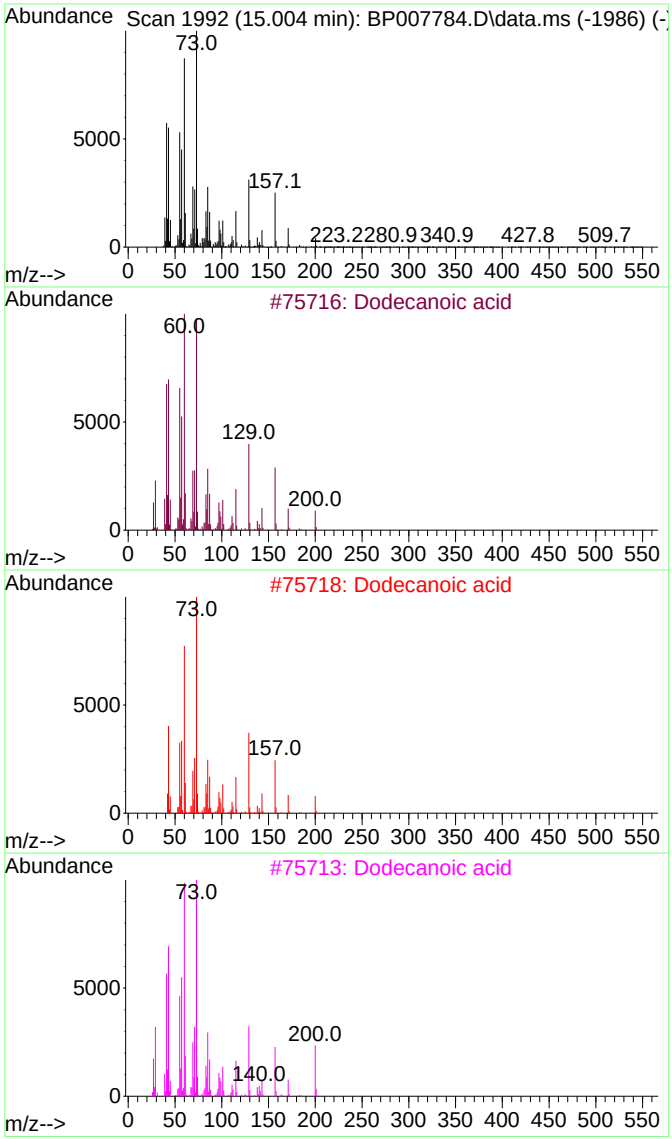
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

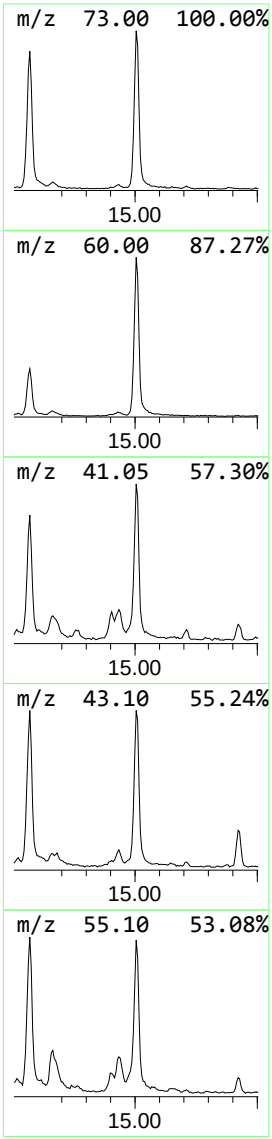
****

Peak Number 20 Dodecanoic acid Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.004	7.63 ng/ul	1355290	Acenaphthene-d10	14.569

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dodecanoic acid	200	C12H24O2	000143-07-7	98
2	Dodecanoic acid	200	C12H24O2	000143-07-7	91
3	Dodecanoic acid	200	C12H24O2	000143-07-7	91
4	Dodecanoic acid	200	C12H24O2	000143-07-7	91
5	Undecanoic acid	186	C11H22O2	000112-37-8	72





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Acq On : 29 Oct 2021 14:44
Operator : CG/JU
Sample : M4262-20DL 2X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BFY53DL

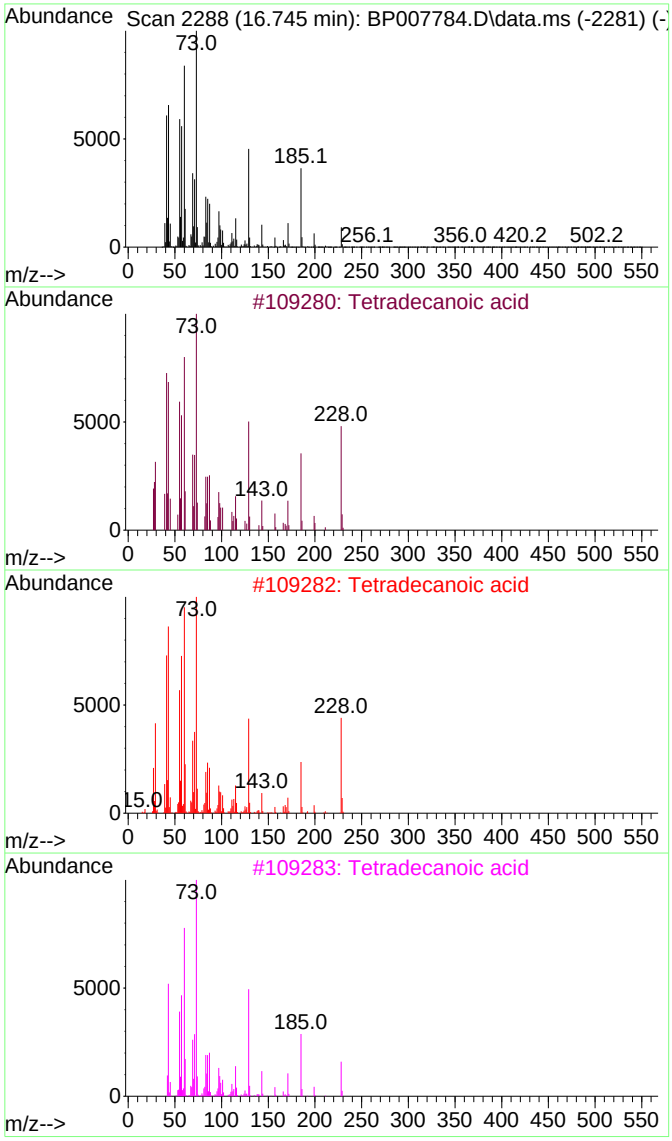
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP102521.M
Quant Title : SVOA CALIBRATION

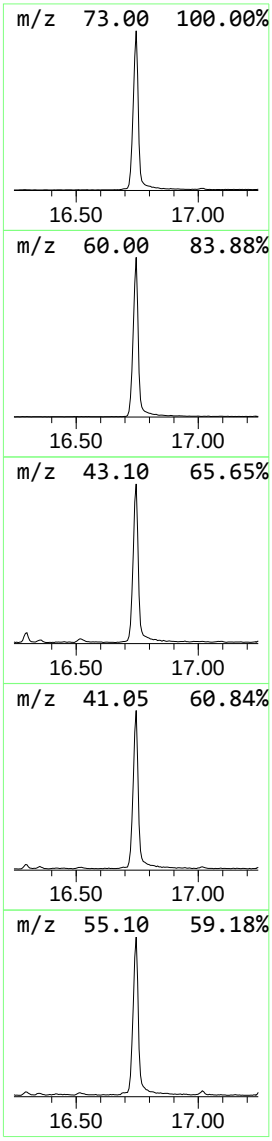
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 21 Tetradecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.745	31.38 ng/ul	4442220	Phenanthrene-d10	17.322

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetradecanoic acid	228	C14H28O2	000544-63-8	96
2	Tetradecanoic acid	228	C14H28O2	000544-63-8	94
3	Tetradecanoic acid	228	C14H28O2	000544-63-8	94
4	Tetradecanoic acid	228	C14H28O2	000544-63-8	91
5	Dodecanoic acid	200	C12H24O2	000143-07-7	68





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Acq On : 29 Oct 2021 14:44
Operator : CG/JU
Sample : M4262-20DL 2X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BFY53DL

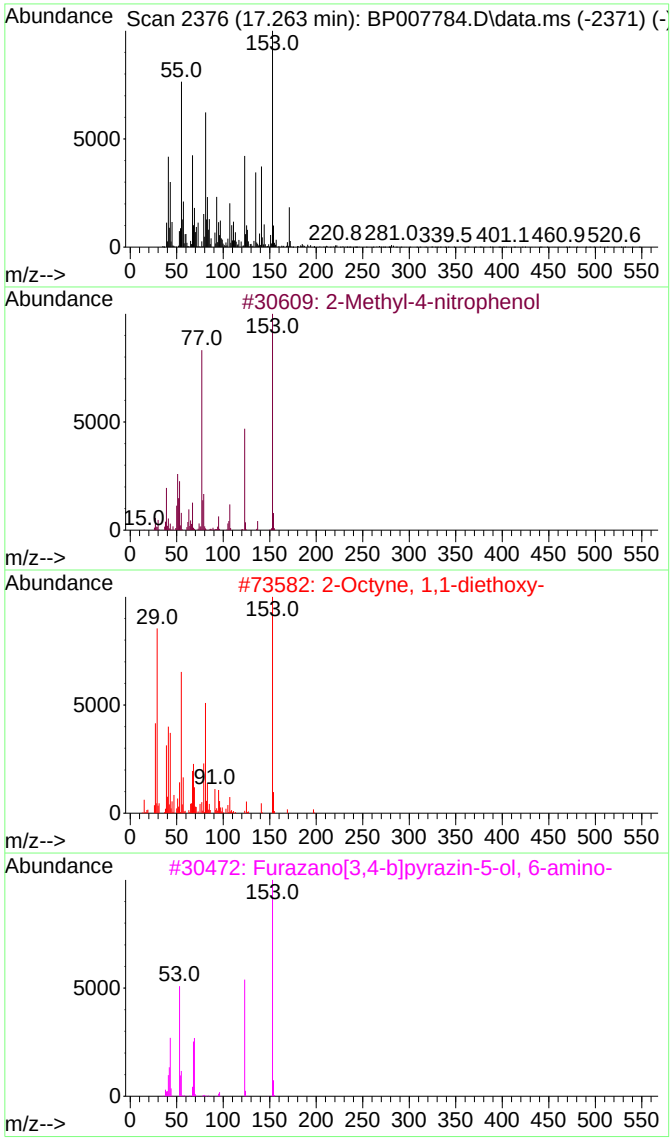
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP102521.M
Quant Title : SVOA CALIBRATION

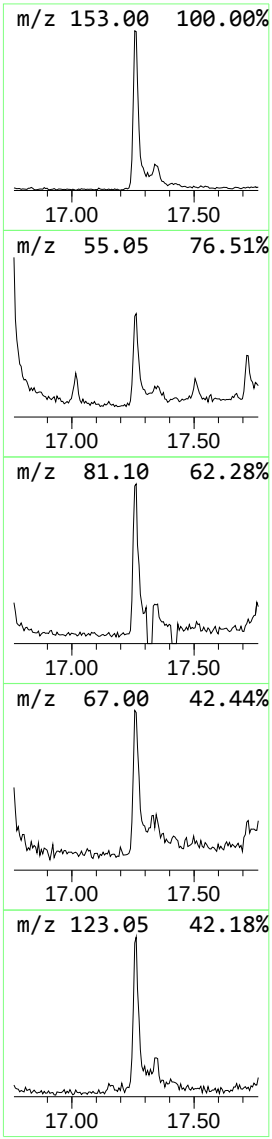
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 22 unknown-04 Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.263	2.01 ng/ul	284365	Phenanthrene-d10	17.322

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Methyl-4-nitrophenol	153	C7H7NO3	000099-53-6	43
2	2-Octyne, 1,1-diethoxy-	198	C12H22O2	016387-55-6	38
3	Furazano[3,4-b]pyrazin-5-ol, 6-a...	153	C4H3N5O2	211919-08-3	38
4	6-Methyl-3-nitro-2-pyridinamine	153	C6H7N3O2	021901-29-1	35
5	4-Nitro-1,3-phenylenediamine	153	C6H7N3O2	005131-58-8	35





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Acq On : 29 Oct 2021 14:44
Operator : CG/JU
Sample : M4262-20DL 2X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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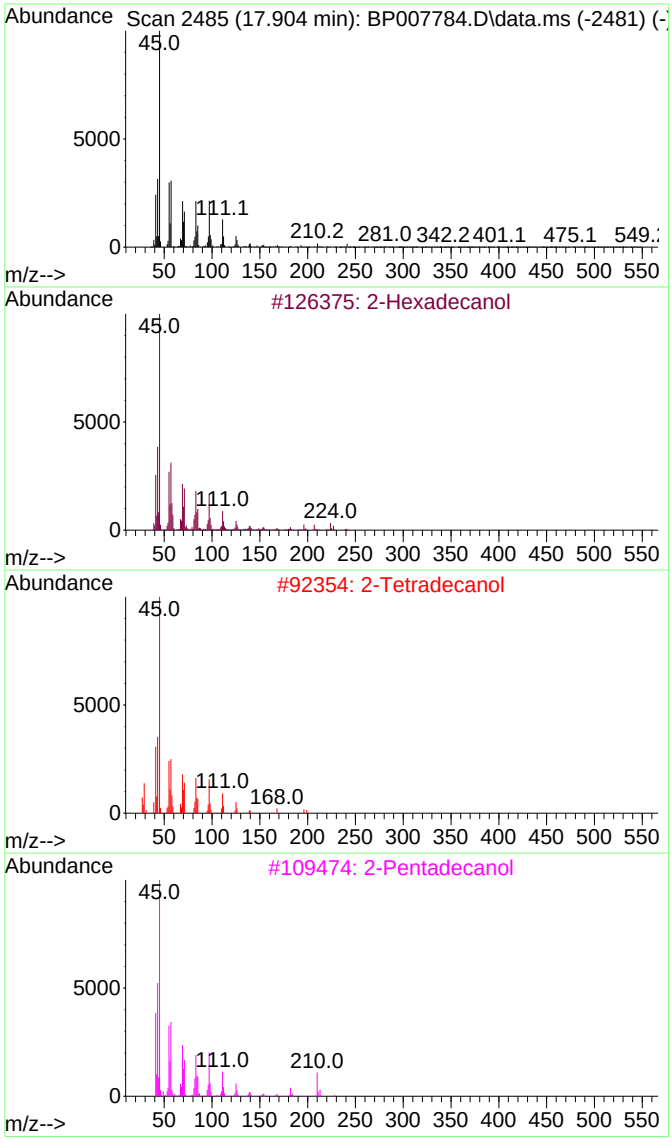
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP102521.M
Quant Title : SVOA CALIBRATION

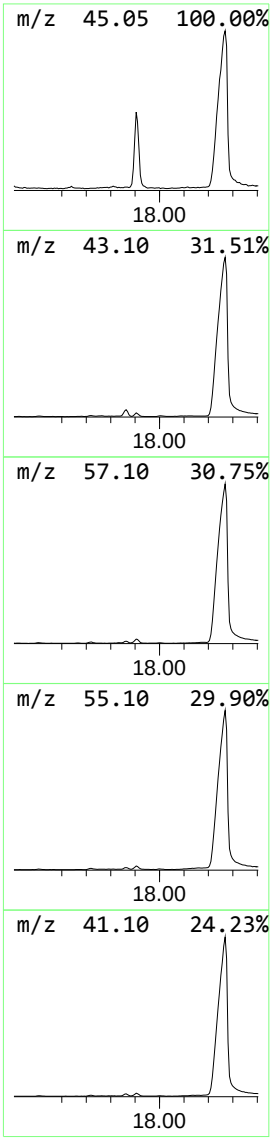
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 23 2-Hexadecanol Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.904	2.53 ng/ul	357602	Phenanthrene-d10	17.322

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Hexadecanol	242	C16H34O	014852-31-4	94
2	2-Tetradecanol	214	C14H30O	004706-81-4	91
3	2-Pentadecanol	228	C15H32O	001653-34-5	90
4	2-Hexadecanol	242	C16H34O	014852-31-4	90
5	2-Heptadecanol	256	C17H36O	016813-18-6	87





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Operator : CG/JU
Sample : M4262-20DL 2X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BFY53DL

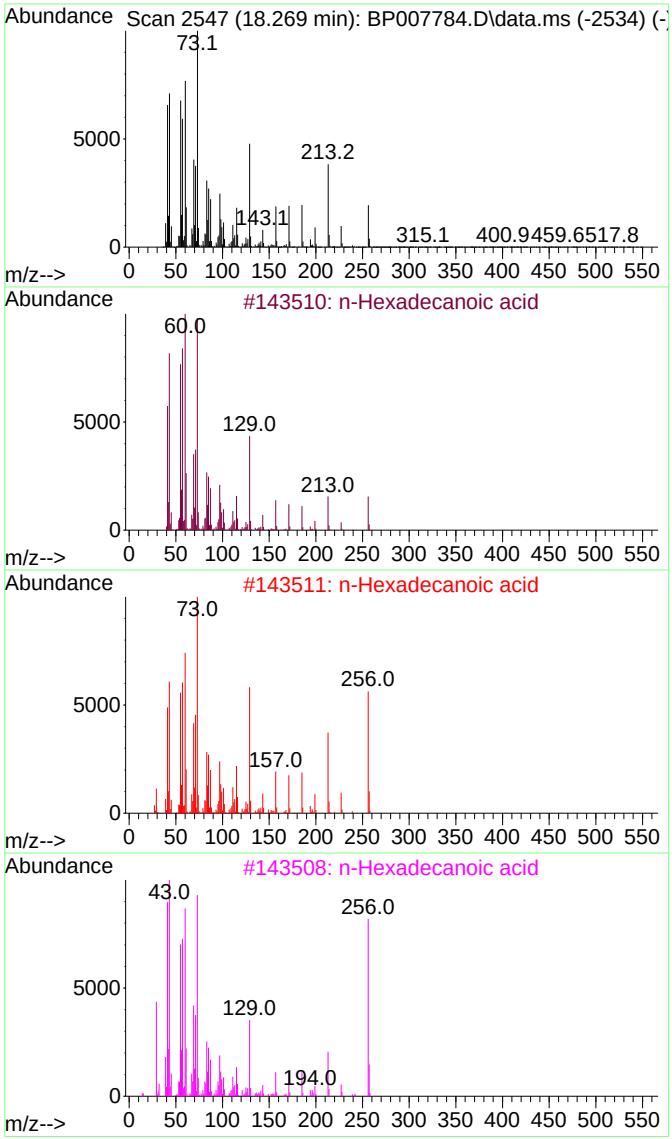
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP102521.M
Quant Title : SVOA CALIBRATION

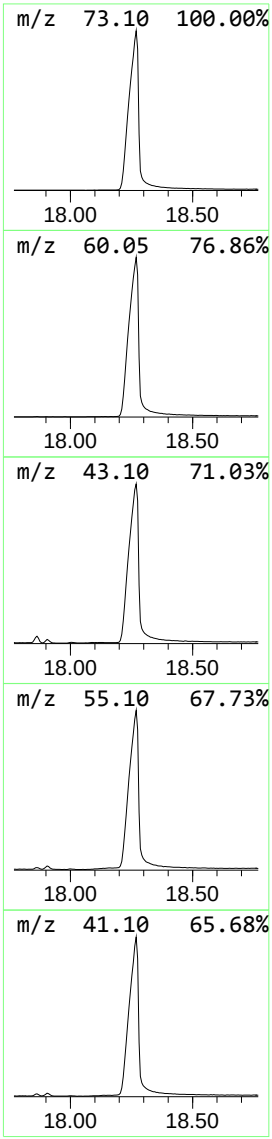
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 24 n-Hexadecanoic acid Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.269	269.70 ng/ul	38175200	Phenanthrene-d10	17.322

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	98
3	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
4	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	95
5	Tridecanoic acid	214	C13H26O2	000638-53-9	93





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Acq On : 29 Oct 2021 14:44
Operator : CG/JU
Sample : M4262-20DL 2X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : BFY53DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP102521.M
Quant Title : SVOA CALIBRATION

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

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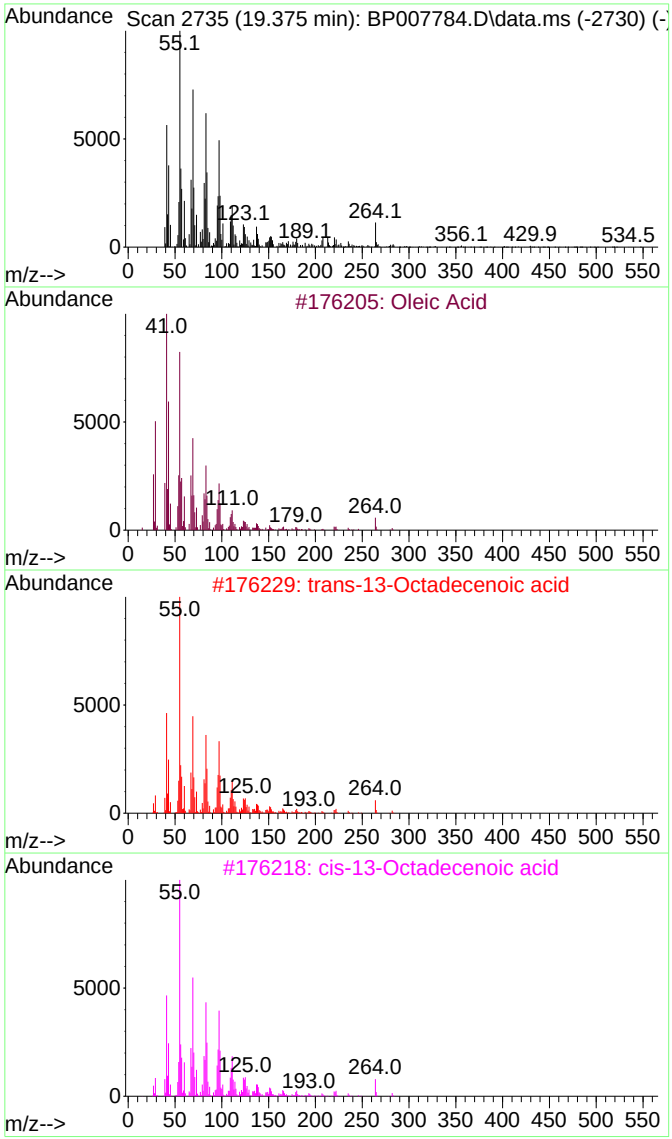
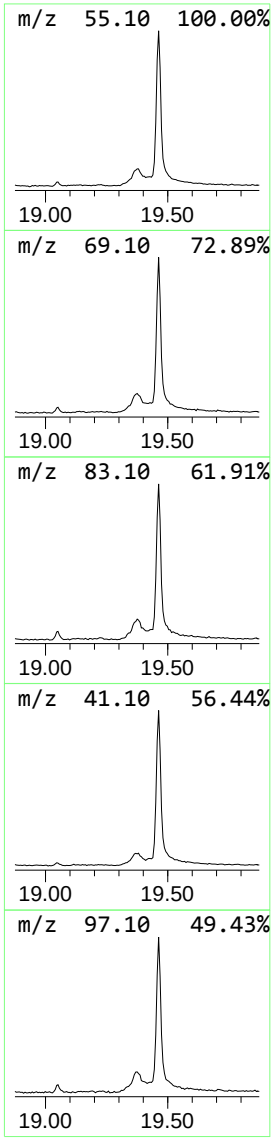
Peak Number 25 Oleic Acid Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.375	2.87 ng/ul	490507	Chrysene-d12	21.410

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Oleic Acid	282	C18H34O2	000112-80-1	99
2	trans-13-Octadecenoic acid	282	C18H34O2	000693-71-0	91
3	cis-13-Octadecenoic acid	282	C18H34O2	013126-39-1	90
4	Palmitoleic acid	254	C16H30O2	000373-49-9	87
5	cis-Vaccenic acid	282	C18H34O2	000506-17-2	87





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Data File : BP007784.D
Acq On : 29 Oct 2021 14:44
Operator : CG/JU
Sample : M4262-20DL 2X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : BFY53DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP102521.M
Quant Title : SVOA CALIBRATION

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

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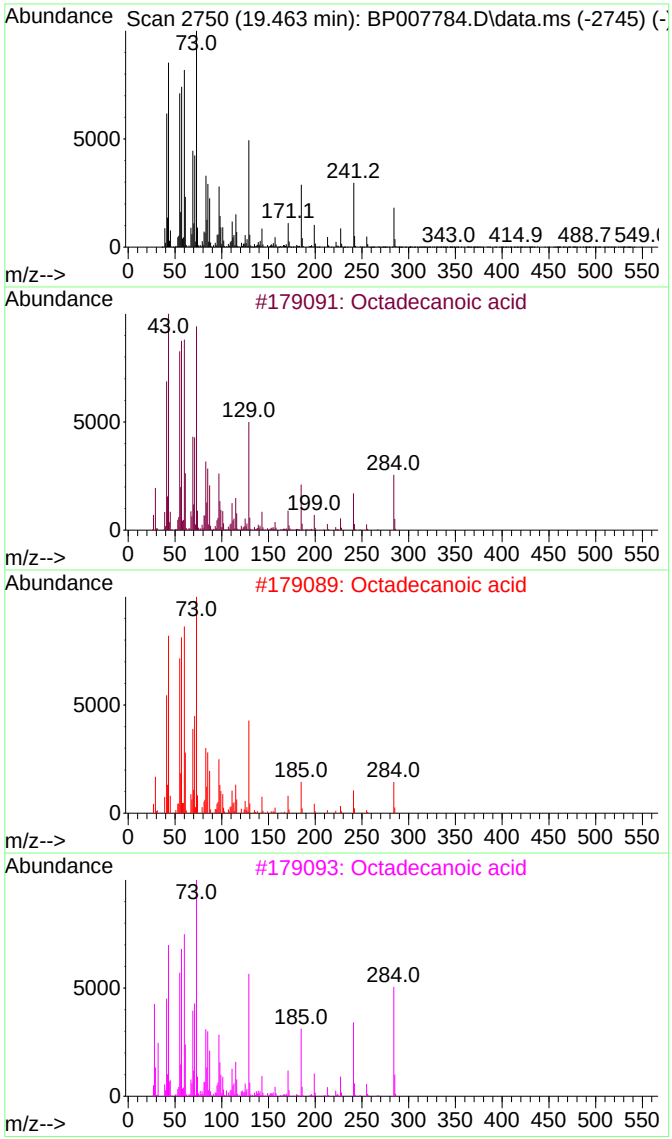
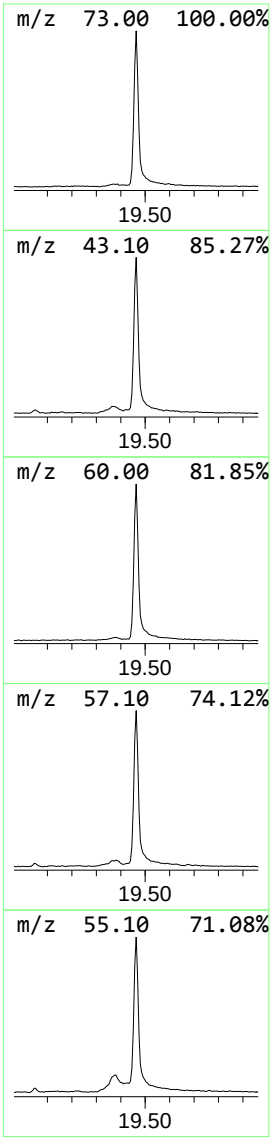
Peak Number 26 Octadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.463	38.03 ng/ul	6502990	Chrysene-d12	21.410

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
4	Octadecanoic acid	284	C18H36O2	000057-11-4	98
5	Octadecanoic acid	284	C18H36O2	000057-11-4	94





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102921\
Data File : BP007784.D
Acq On : 29 Oct 2021 14:44
Operator : CG/JU
Sample : M4262-20DL 2X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : BFY53DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP102521.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,...	3.681	2.3	ng/ul	184201	1	7.928	1575920	20.0
Butanoic acid	4.111	25.5	ng/ul	2006800	1	7.928	1575920	20.0
Butanoic acid, ...	4.834	5.7	ng/ul	450762	1	7.928	1575920	20.0
Butanoic acid, ...	5.058	22.1	ng/ul	1742410	1	7.928	1575920	20.0
Pentanoic acid	5.534	28.3	ng/ul	2232030	1	7.928	1575920	20.0
5-Methylhexanoi...	8.016	2.2	ng/ul	171408	1	7.928	1575920	20.0
Hexanoic acid, ...	8.181	3.7	ng/ul	294880	1	7.928	1575920	20.0
unknown-01	9.557	5.3	ng/ul	414612	2	10.728	1558920	20.0
unknown-02	9.599	2.5	ng/ul	191975	2	10.728	1558920	20.0
Ethanol, 2-(2-b...	10.522	6.0	ng/ul	464508	2	10.728	1558920	20.0
Nonanoic acid	11.587	14.9	ng/ul	1157370	2	10.728	1558920	20.0
Decanoic acid, ...	12.222	2.6	ng/ul	205811	2	10.728	1558920	20.0
Decanoic acid, ...	12.322	9.9	ng/ul	772335	2	10.728	1558920	20.0
(DEL) Alkane: S...	12.399	3.9	ng/ul	300758	2	10.728	1558920	20.0
unknown-03	12.451	4.9	ng/ul	378701	2	10.728	1558920	20.0
9-Decenoic acid	12.828	3.9	ng/ul	684636	3	14.569	3550580	20.0
n-Decanoic acid	12.940	50.7	ng/ul	9000110	3	14.569	3550580	20.0
Benzenebutanoic...	13.722	4.6	ng/ul	815609	3	14.569	3550580	20.0
9-Tetradecenal,...	14.663	2.4	ng/ul	426703	3	14.569	3550580	20.0
Dodecanoic acid	15.004	7.6	ng/ul	1355290	3	14.569	3550580	20.0
Tetradecanoic acid	16.745	31.4	ng/ul	4442220	4	17.322	2830920	20.0
unknown-04	17.263	2.0	ng/ul	284365	4	17.322	2830920	20.0
2-Hexadecanol	17.904	2.5	ng/ul	357602	4	17.322	2830920	20.0
n-Hexadecanoic ...	18.269	269.7	ng/ul	38175200	4	17.322	2830920	20.0
Oleic Acid	19.375	2.9	ng/ul	490507	5	21.410	3419840	20.0
Octadecanoic acid	19.463	38.0	ng/ul	6502990	5	21.410	3419840	20.0