

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102121\
 Data File : BP007603.D
 Acq On : 22 Oct 2021 23:10
 Operator : CG/JU
 Sample : M4281-05DL 5X
 Misc :
 ALS Vial : 63 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BFY61DL

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: BP007603.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.376	12	15	17	rBV	12384	11795	0.33%	0.034%
2	3.411	17	21	30	rVB	133548	174321	4.89%	0.507%
3	3.670	58	65	72	rBV	52799	96676	2.71%	0.281%
4	3.805	86	88	98	rBV	48904	83064	2.33%	0.242%
5	4.052	115	130	137	rBV	219266	595318	16.70%	1.732%
6	4.123	139	142	148	rVB2	12416	18444	0.52%	0.054%
7	4.817	250	260	272	rBV2	85141	254426	7.14%	0.740%
8	5.017	275	294	304	rBV	281023	946336	26.55%	2.753%
9	5.517	353	379	382	rBV	428194	1848492	51.85%	5.378%
10	5.834	428	433	438	rVB9	5364	8645	0.24%	0.025%
11	5.981	452	458	463	rBV10	3346	8503	0.24%	0.025%
12	6.028	463	466	471	rVB5	3718	5431	0.15%	0.016%
13	6.405	510	530	540	rBV	319942	1037683	29.11%	3.019%
14	6.481	540	543	546	rVB5	12811	15898	0.45%	0.046%
15	6.811	595	599	603	rVB5	3930	5141	0.14%	0.015%
16	7.034	617	637	645	rBV	425704	1481469	41.56%	4.310%
17	7.105	645	649	659	rVB	48733	92695	2.60%	0.270%
18	7.269	671	677	687	rVB	137666	227287	6.38%	0.661%
19	7.475	705	712	718	rBV	143378	223168	6.26%	0.649%
20	7.552	723	725	732	rVB6	3870	6359	0.18%	0.019%
21	7.911	775	786	788	rBV	209158	459685	12.90%	1.337%
22	7.946	788	792	799	rVB	754226	1163318	32.63%	3.384%
23	8.146	823	826	830	rVB6	4628	6356	0.18%	0.018%
24	8.199	830	835	838	rBV7	3205	5686	0.16%	0.017%
25	8.469	876	881	885	rBV3	6052	9986	0.28%	0.029%
26	8.563	885	897	906	rVV	273067	634435	17.80%	1.846%
27	8.652	906	912	917	rVV	123249	209858	5.89%	0.611%
28	8.705	917	921	927	rVB	25937	43024	1.21%	0.125%
29	8.840	939	944	945	rBV5	2930	4141	0.12%	0.012%
30	9.099	980	988	998	rBV	170246	283009	7.94%	0.823%
31	9.222	1007	1009	1013	rBV5	2949	4778	0.13%	0.014%
32	9.410	1035	1041	1052	rBV	43019	79809	2.24%	0.232%
33	9.563	1063	1067	1073	rVB7	7578	13405	0.38%	0.039%
34	9.716	1090	1093	1096	rVB5	2457	3708	0.10%	0.011%
35	9.828	1105	1112	1120	rBV	114906	195051	5.47%	0.567%
36	10.105	1146	1159	1173	rBV	386843	889458	24.95%	2.588%

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Integrator: RTE

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Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP101421.M
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37	10.369	1197	1204	1214	rVB	211827	367110	10.30%	1.068%
38	10.534	1225	1232	1239	rBV	158935	262651	7.37%	0.764%
39	10.581	1239	1240	1243	rVB3	4079	3653	0.10%	0.011%
40	10.746	1261	1268	1278	rVB	994348	1725214	48.40%	5.019%
41	10.881	1284	1291	1299	rBV	166226	286061	8.02%	0.832%
42	11.075	1320	1324	1326	rBV5	3117	3922	0.11%	0.011%
43	11.157	1335	1338	1344	rVB7	6580	11164	0.31%	0.032%
44	11.275	1353	1358	1365	rVV	13286	24810	0.70%	0.072%
45	11.363	1369	1373	1375	rVB5	2496	3575	0.10%	0.010%
46	11.393	1375	1378	1386	rBV10	4602	8869	0.25%	0.026%
47	11.546	1397	1404	1416	rBV2	32621	64249	1.80%	0.187%
48	12.063	1488	1492	1498	rVB6	10850	17417	0.49%	0.051%
49	12.334	1531	1538	1540	rBV6	6158	10958	0.31%	0.032%
50	12.363	1540	1543	1548	rVB7	3419	6959	0.20%	0.020%
51	12.540	1569	1573	1580	rBV4	10283	16995	0.48%	0.049%
52	12.722	1600	1604	1610	rVB8	5337	10526	0.30%	0.031%
53	12.781	1610	1614	1617	rBV6	2598	4033	0.11%	0.012%
54	12.869	1623	1629	1639	rBV	121234	180412	5.06%	0.525%
55	13.204	1681	1686	1690	rVB7	5053	8688	0.24%	0.025%
56	13.322	1701	1706	1708	rBV6	3497	5462	0.15%	0.016%
57	13.493	1729	1735	1740	rBV3	14450	22891	0.64%	0.067%
58	13.551	1740	1745	1746	rBV5	3649	4581	0.13%	0.013%
59	13.710	1769	1772	1776	rBV6	3663	4757	0.13%	0.014%
60	13.987	1813	1819	1825	rBV	439341	608815	17.08%	1.771%
61	14.269	1858	1867	1874	rBV	457382	695703	19.52%	2.024%
62	14.428	1890	1894	1896	rBV5	3250	3912	0.11%	0.011%
63	14.475	1896	1902	1903	rVV6	6393	9279	0.26%	0.027%
64	14.504	1903	1907	1910	rVV3	8306	15269	0.43%	0.044%
65	14.581	1913	1920	1931	rVV	1640296	2435963	68.33%	7.087%
66	14.769	1947	1952	1958	rVB2	32008	47103	1.32%	0.137%
67	14.845	1963	1965	1969	rVB5	3116	3797	0.11%	0.011%
68	15.004	1987	1992	1993	rBV5	7179	9988	0.28%	0.029%
69	15.034	1993	1997	2002	rVB8	8296	16420	0.46%	0.048%
70	15.134	2007	2014	2038	rBV3	71841	400114	11.22%	1.164%
71	15.322	2044	2046	2052	rVB5	10044	14060	0.39%	0.041%
72	15.440	2063	2066	2070	rVB4	8403	11928	0.33%	0.035%
73	15.575	2082	2089	2097	rBV	668885	970240	27.22%	2.823%
74	15.687	2103	2108	2113	rVB4	20102	29493	0.83%	0.086%
75	15.816	2126	2130	2135	rVB4	9521	14554	0.41%	0.042%
76	15.934	2149	2150	2154	rBV4	3828	5014	0.14%	0.015%

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Integration Parameters: LSCINT.P

Integrator: RTE
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 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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 Title : SVOA CALIBRATION

77	16.016	2162	2164	2169	rVB6	5408	8155	0.23%	0.024%
78	16.134	2180	2184	2186	rBV4	4280	5470	0.15%	0.016%
79	16.310	2212	2214	2218	rVB5	4624	4613	0.13%	0.013%
80	16.369	2222	2224	2228	rVV4	3851	4143	0.12%	0.012%
81	16.528	2249	2251	2255	rBV5	6038	6672	0.19%	0.019%
82	16.992	2326	2330	2337	rBV2	8892	22015	0.62%	0.064%
83	17.081	2343	2345	2348	rBV3	9952	13810	0.39%	0.040%
84	17.334	2381	2388	2397	rBV	2027012	2969443	83.30%	8.639%
85	17.434	2399	2405	2412	rVV	792362	1126035	31.59%	3.276%
86	18.016	2500	2504	2509	rVB	46045	51644	1.45%	0.150%
87	18.228	2536	2540	2547	rBV2	90429	121076	3.40%	0.352%
88	18.522	2587	2590	2596	rBV5	22381	37292	1.05%	0.108%
89	19.145	2693	2696	2700	rBV3	48105	57258	1.61%	0.167%
90	19.610	2771	2775	2779	rBV2	83794	105528	2.96%	0.307%
91	19.663	2779	2784	2790	rVV	1059367	1404885	39.41%	4.087%
92	20.622	2945	2947	2952	rVB3	67907	73240	2.05%	0.213%
93	21.133	3031	3034	3040	rBV5	104805	153524	4.31%	0.447%
94	21.416	3077	3082	3088	rBV	2531079	3423202	96.03%	9.959%
95	23.704	3465	3471	3479	rVB2	675344	1365508	38.31%	3.973%
96	23.792	3483	3486	3491	rVV	219029	360310	10.11%	1.048%
97	23.863	3492	3498	3510	rVB2	1662907	3564785	100.00%	10.371%

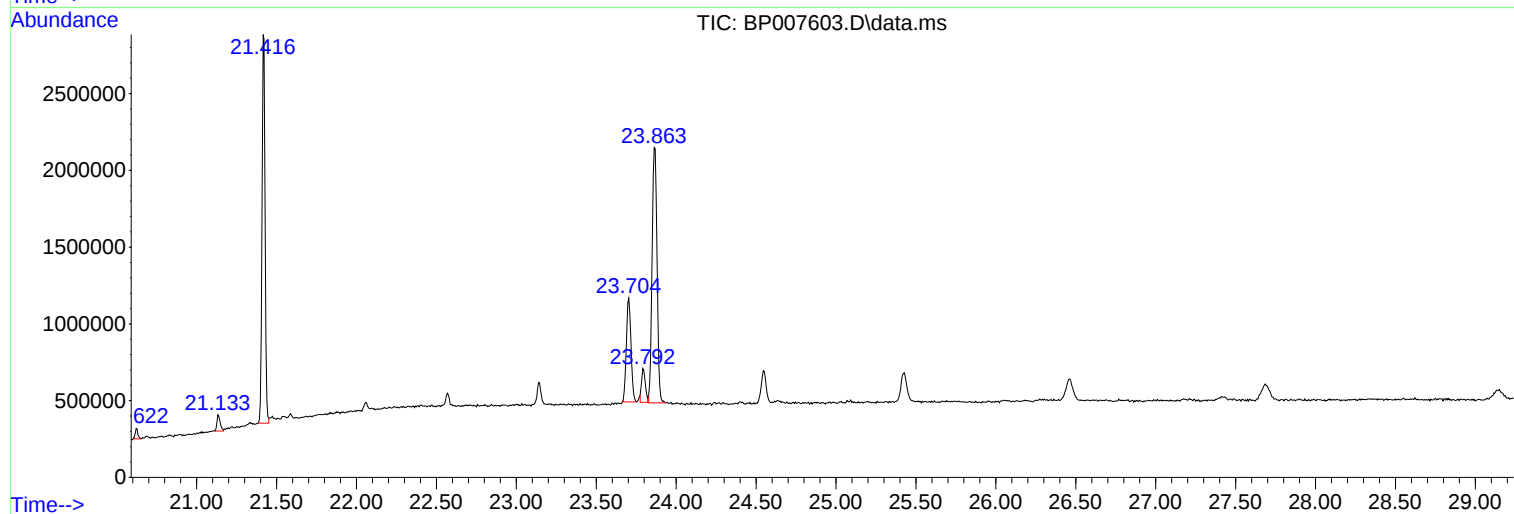
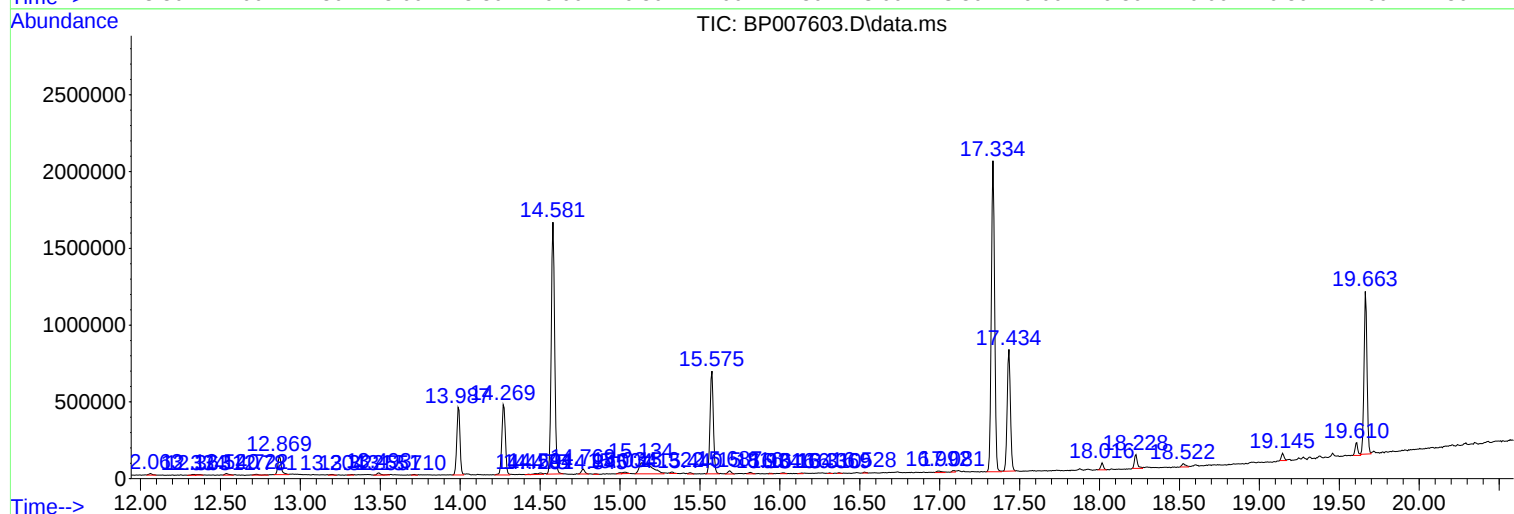
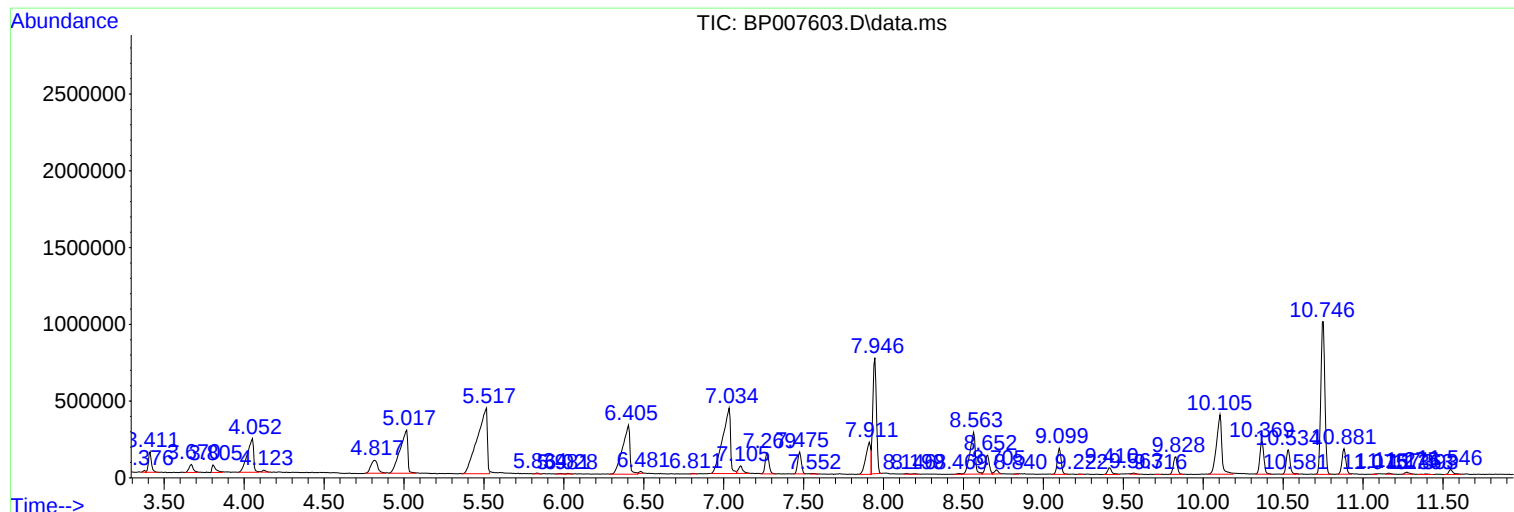
Sum of corrected areas: 34372075

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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



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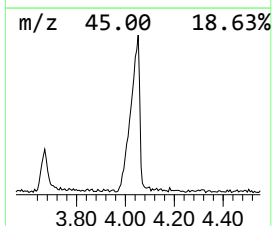
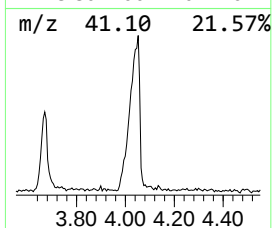
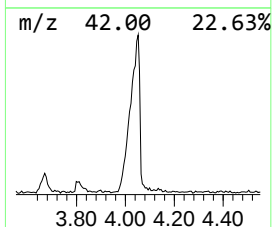
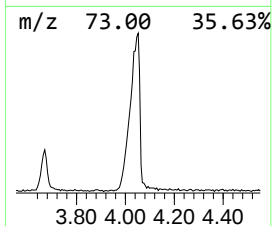
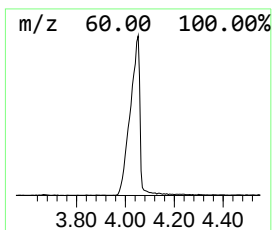
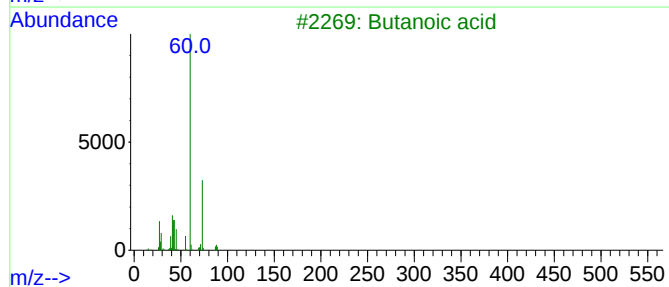
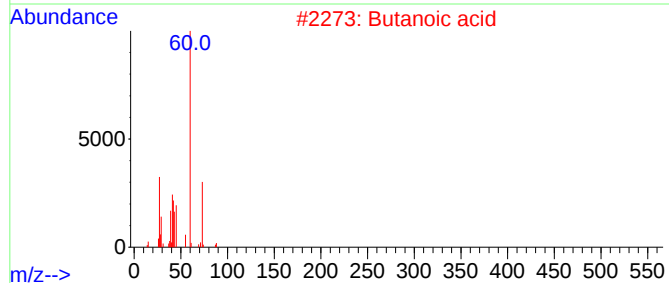
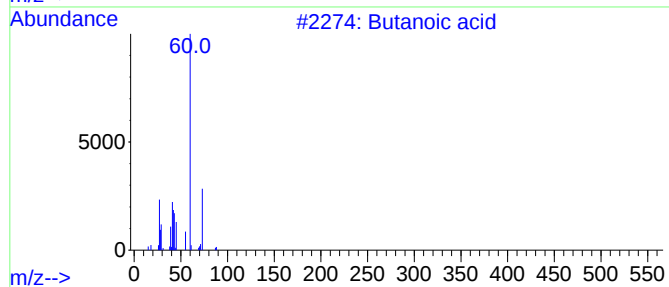
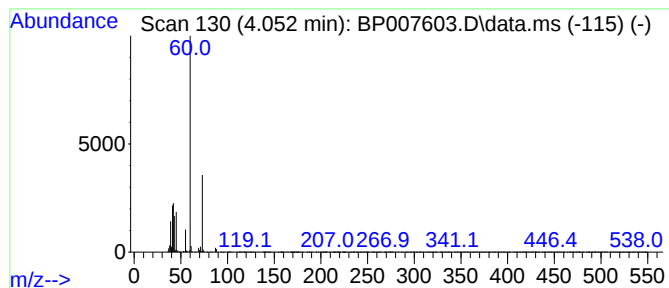
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP101421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Butanoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.052	10.23 ng/ul	595318	1,4-Dichlorobenzene-d4	7.946

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	90
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	72



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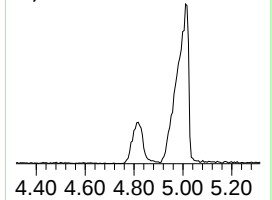
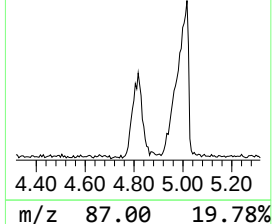
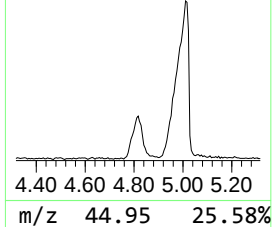
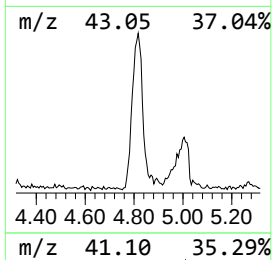
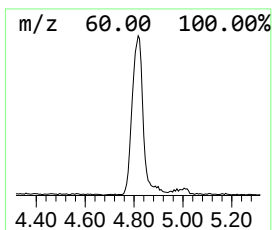
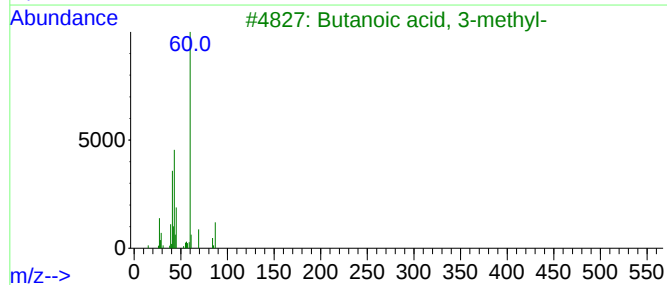
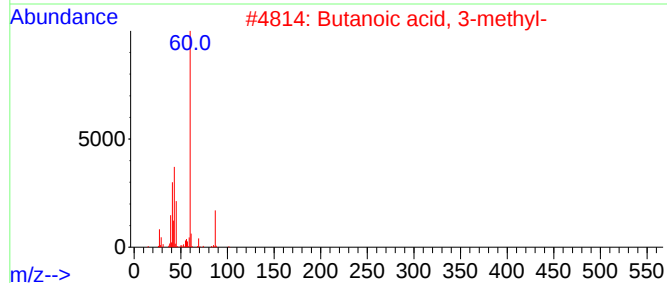
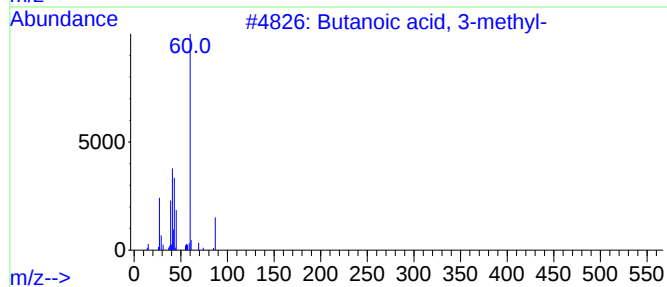
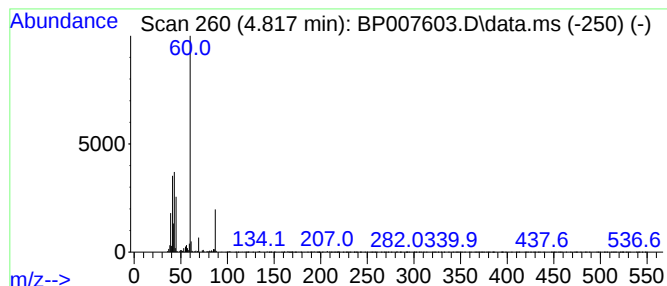
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP101421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Butanoic acid, 3-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.817	4.37 ng/ul	254426	1,4-Dichlorobenzene-d4	7.946

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	59
4			Butanoic acid, 3-methyl-	102	C5H10O2	000503-74-2	46
5			Butanoic acid, 3-methyl-	102	C5H10O2	000503-74-2	38



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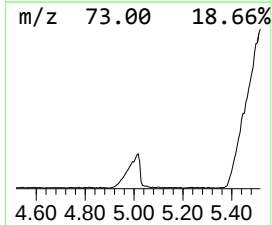
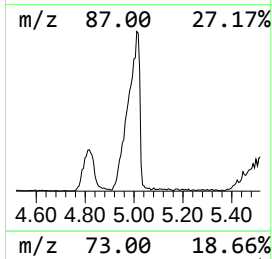
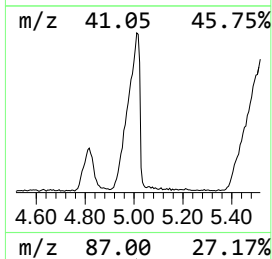
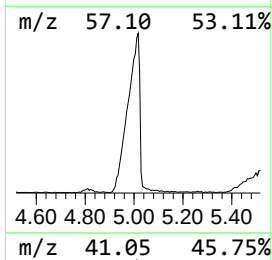
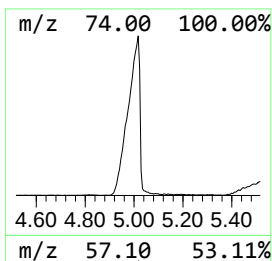
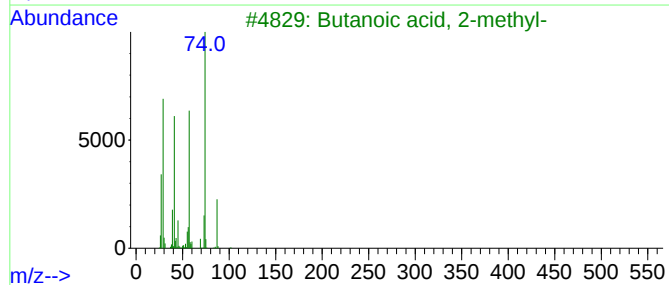
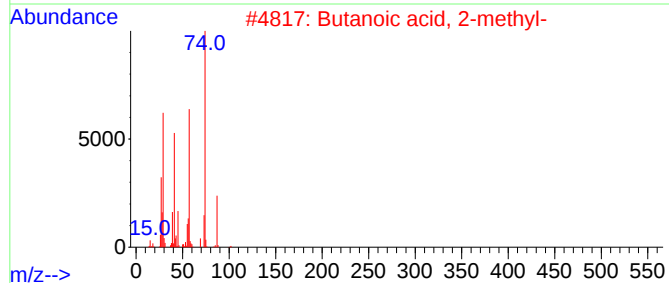
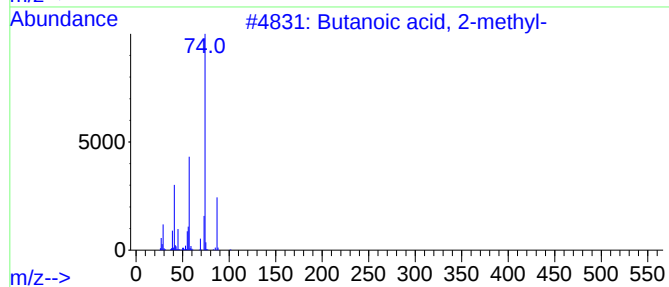
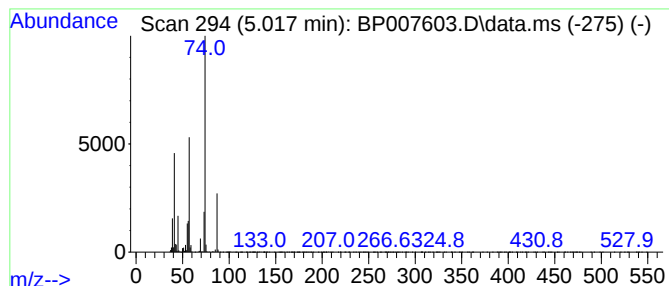
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 Butanoic acid, 2-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.017	16.27 ng/ul	946336	1,4-Dichlorobenzene-d4	7.946

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	83
3			Butanoic acid, 2-methyl-	102	C5H10O2	000116-53-0	78
4			Pentanoic acid, 2-methyl-	116	C6H12O2	000097-61-0	56
5			Hexanoic acid, 2-methyl-	130	C7H14O2	004536-23-6	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102121\
Data File : BP007603.D
Acq On : 22 Oct 2021 23:10
Operator : CG/JU
Sample : M4281-05DL 5X
Misc :
ALS Vial : 63 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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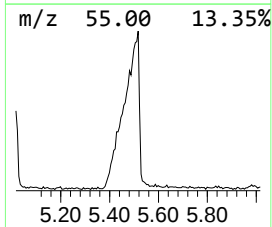
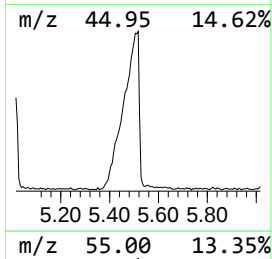
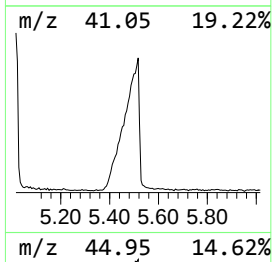
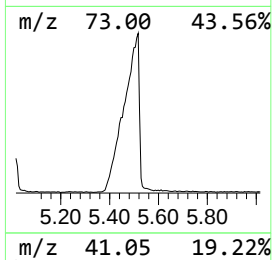
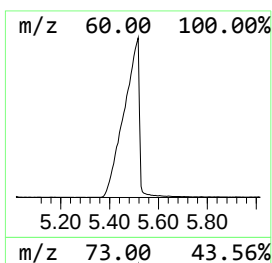
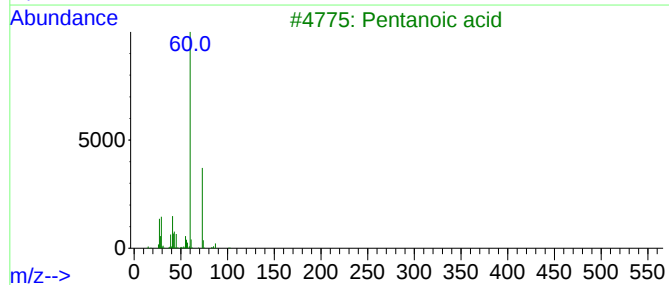
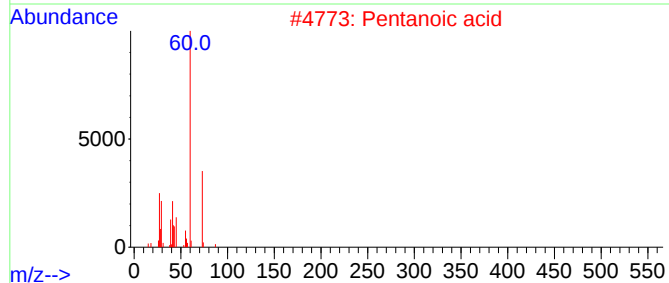
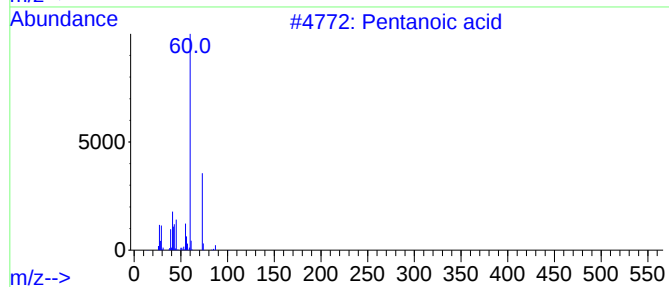
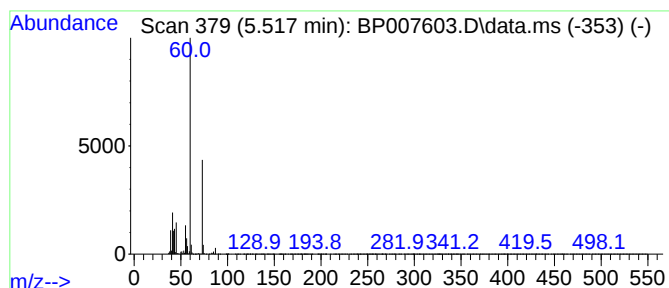
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Pentanoic acid Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.517	31.78 ng/ul	1848490	1,4-Dichlorobenzene-d4	7.946

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			Butanoic acid	88	C4H8O2	000107-92-6	78
5			Pentanoic acid	102	C5H10O2	000109-52-4	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102121\
Data File : BP007603.D
Acq On : 22 Oct 2021 23:10
Operator : CG/JU
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Misc :
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ClientSampleId :
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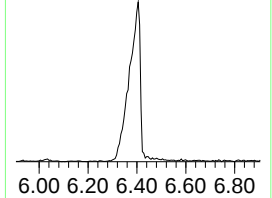
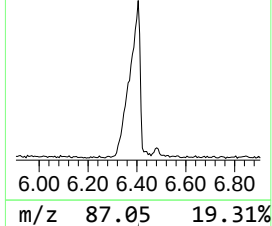
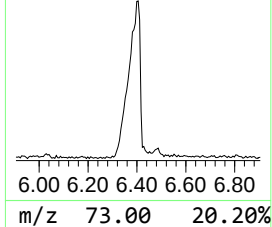
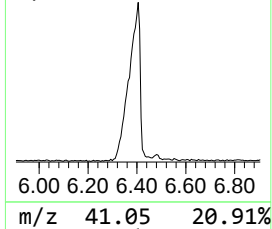
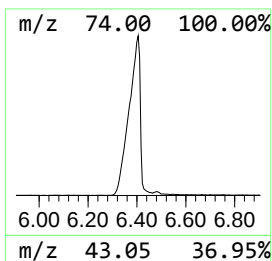
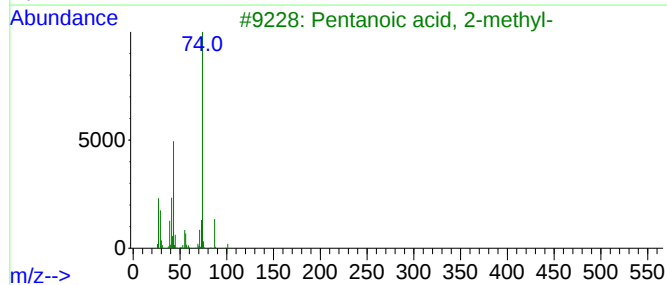
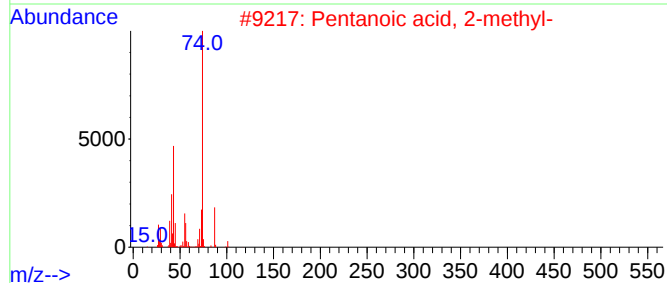
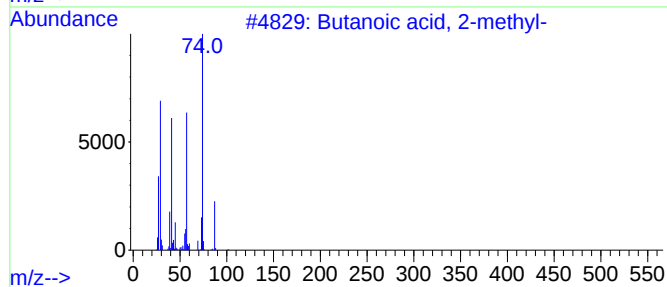
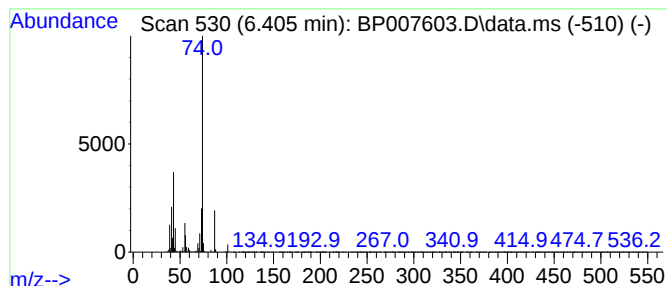
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Pentanoic acid, 2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.405	17.84 ng/ul	1037680	1,4-Dichlorobenzene-d4	7.946

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butanoic acid, 2-methyl-	102	C5H10O2	000116-53-0	80
2			Pentanoic acid, 2-methyl-	116	C6H12O2	000097-61-0	78
3			Pentanoic acid, 2-methyl-	116	C6H12O2	000097-61-0	72
4			Pentanoic acid, 2-methyl-	116	C6H12O2	000097-61-0	72
5			Hexanoic acid, 2-methyl-	130	C7H14O2	004536-23-6	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102121\
Data File : BP007603.D
Acq On : 22 Oct 2021 23:10
Operator : CG/JU
Sample : M4281-05DL 5X
Misc :
ALS Vial : 63 Sample Multiplier: 1

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ClientSampleId :
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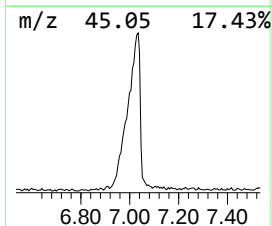
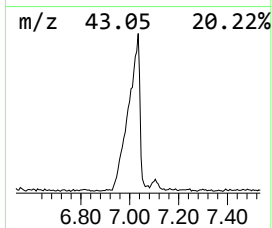
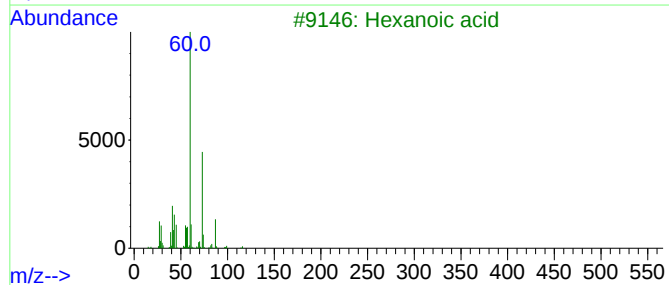
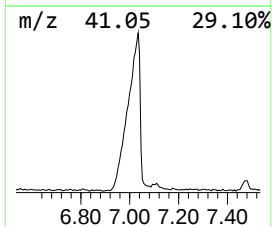
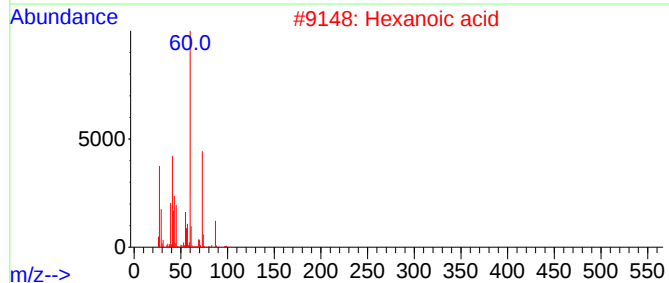
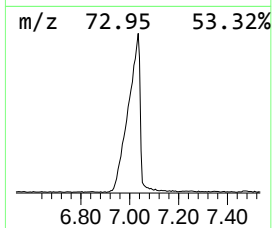
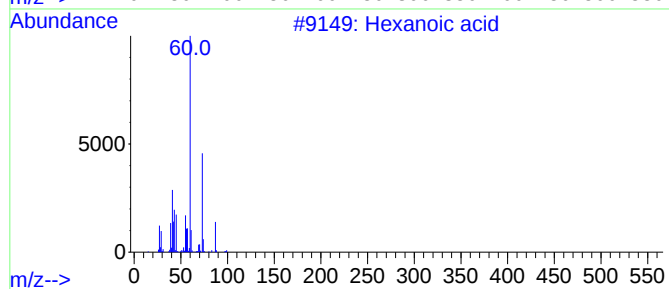
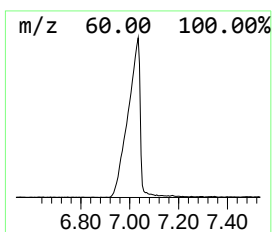
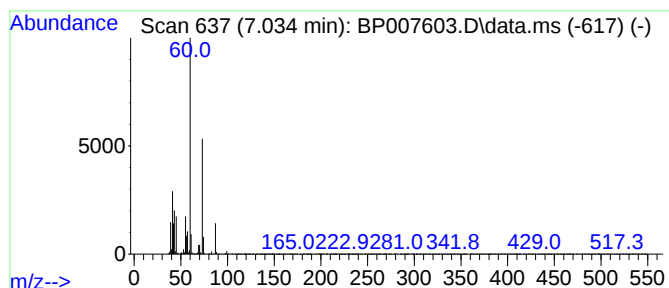
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Hexanoic acid Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.034	25.47 ng/ul	1481470	1,4-Dichlorobenzene-d4	7.946

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	90
3			Hexanoic acid	116	C6H12O2	000142-62-1	83
4			Heptanoic acid	130	C7H14O2	000111-14-8	74
5			Pentanoic acid	102	C5H10O2	000109-52-4	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102121\
Data File : BP007603.D
Acq On : 22 Oct 2021 23:10
Operator : CG/JU
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Misc :
ALS Vial : 63 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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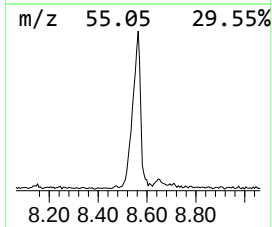
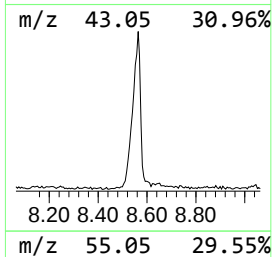
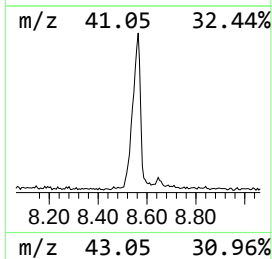
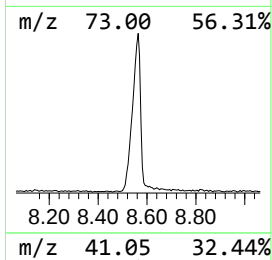
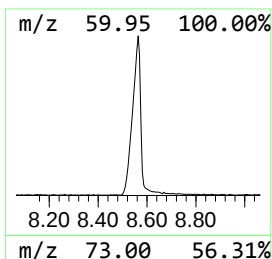
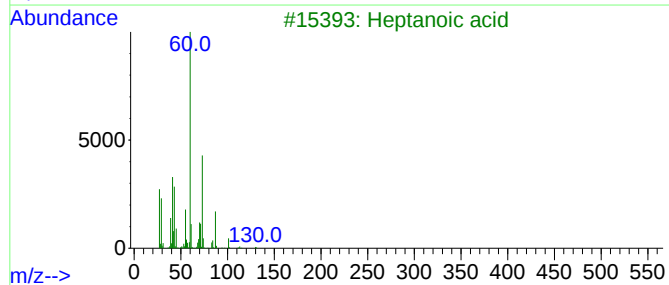
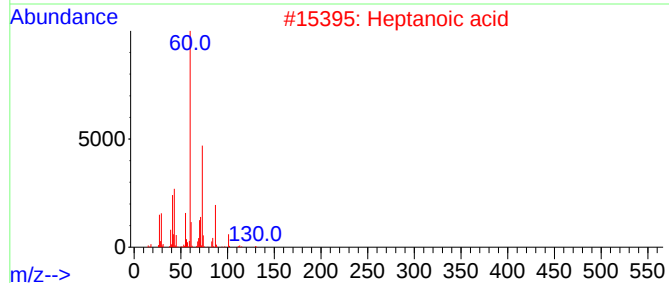
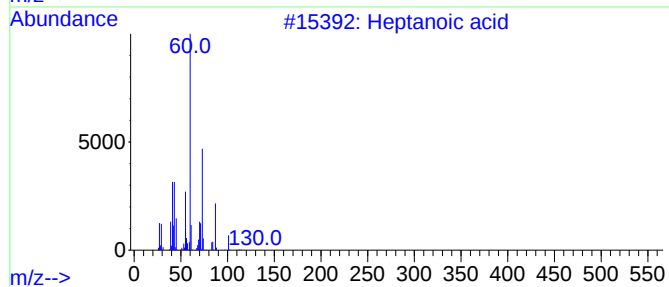
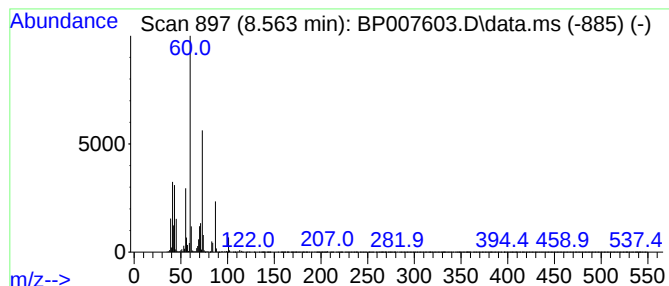
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Heptanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.563	10.91 ng/ul	634435	1,4-Dichlorobenzene-d4	7.946

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptanoic acid	130	C7H14O2	000111-14-8	91
2			Heptanoic acid	130	C7H14O2	000111-14-8	90
3			Heptanoic acid	130	C7H14O2	000111-14-8	90
4			Heptanoic acid	130	C7H14O2	000111-14-8	78
5			Hexanoic acid	116	C6H12O2	000142-62-1	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102121\
Data File : BP007603.D
Acq On : 22 Oct 2021 23:10
Operator : CG/JU
Sample : M4281-05DL 5X
Misc :
ALS Vial : 63 Sample Multiplier: 1

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ClientSampleId :
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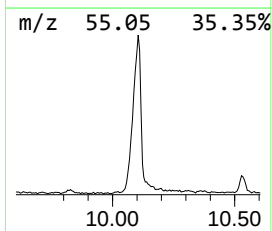
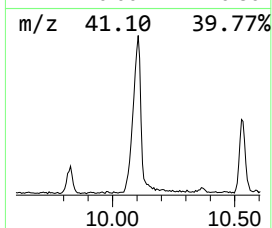
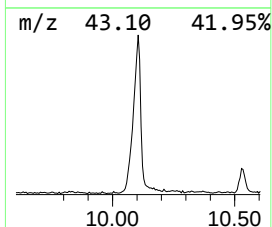
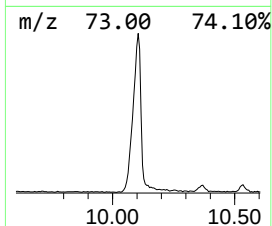
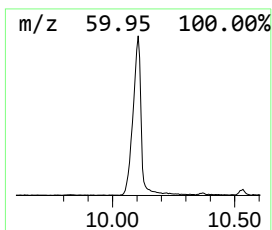
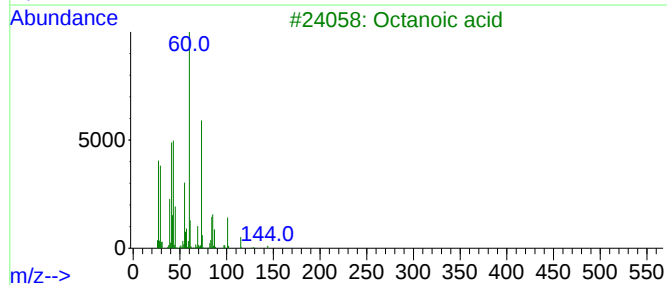
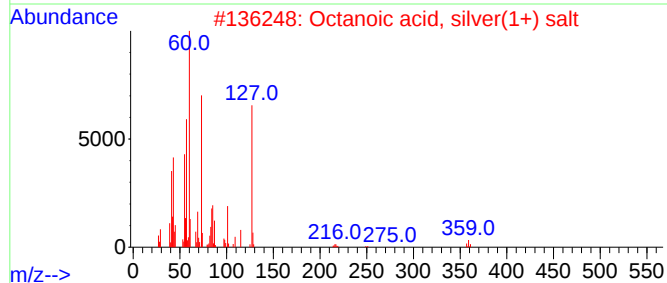
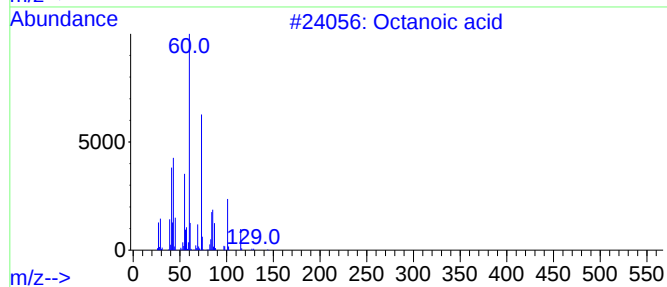
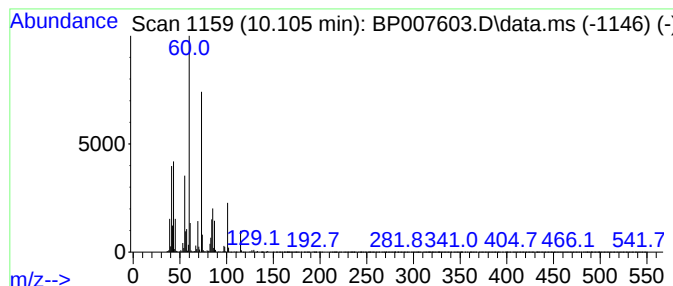
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Octanoic acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.105	10.31 ng/ul	889458	Naphthalene-d8	10.752

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octanoic acid	144	C8H16O2	000124-07-2	97
2			Octanoic acid, silver(1+) salt	250	C8H15AgO2	024927-67-1	87
3			Octanoic acid	144	C8H16O2	000124-07-2	83
4			Octanoic acid	144	C8H16O2	000124-07-2	72
5			Hexanoic acid	116	C6H12O2	000142-62-1	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102121\
Data File : BP007603.D
Acq On : 22 Oct 2021 23:10
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ClientSampleId :
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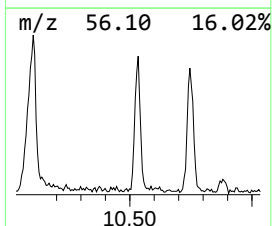
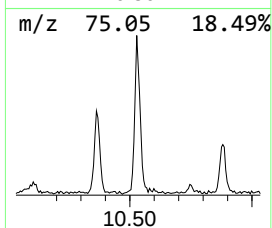
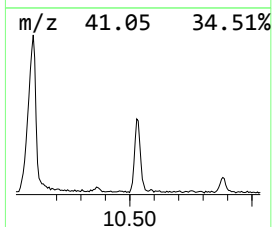
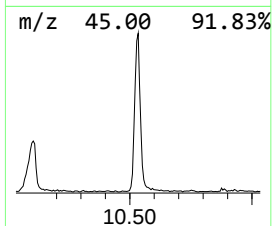
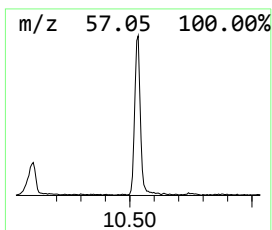
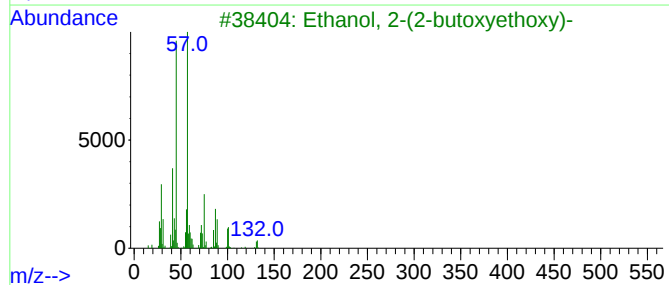
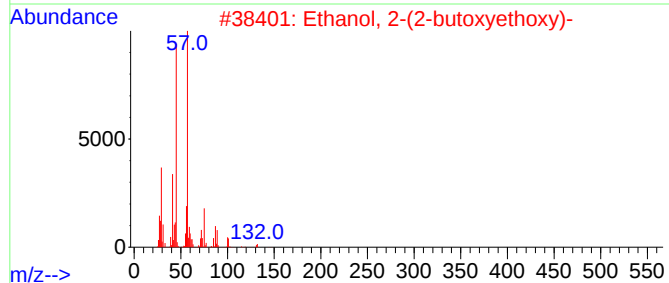
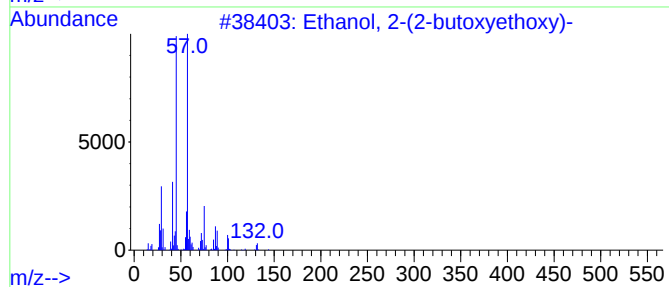
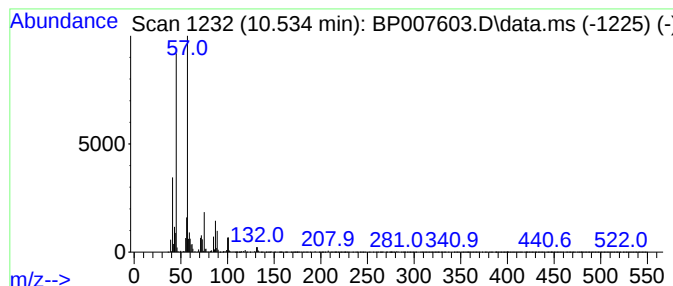
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 Ethanol, 2-(2-butoxyethoxy)- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.534	3.04 ng/ul	262651	Naphthalene-d8	10.752

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	86
4			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	86
5			Ethanol, 1-(2-butoxyethoxy)-	162	C8H18O3	054446-78-5	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102121\
Data File : BP007603.D
Acq On : 22 Oct 2021 23:10
Operator : CG/JU
Sample : M4281-05DL 5X
Misc :
ALS Vial : 63 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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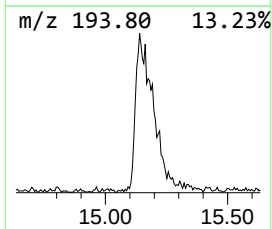
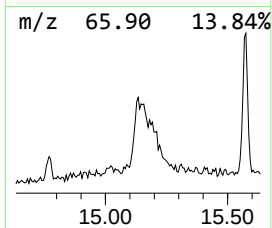
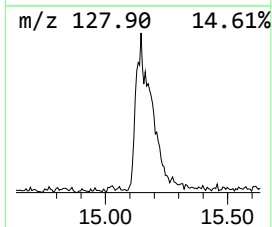
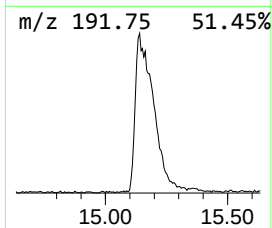
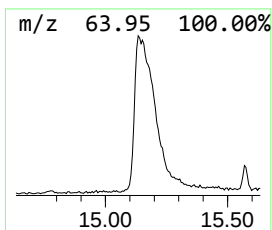
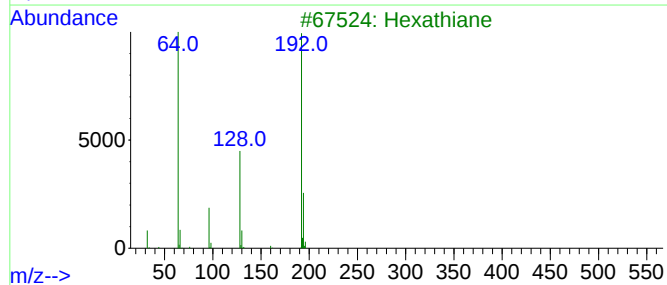
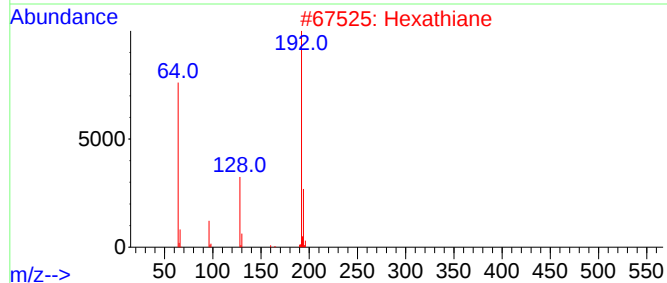
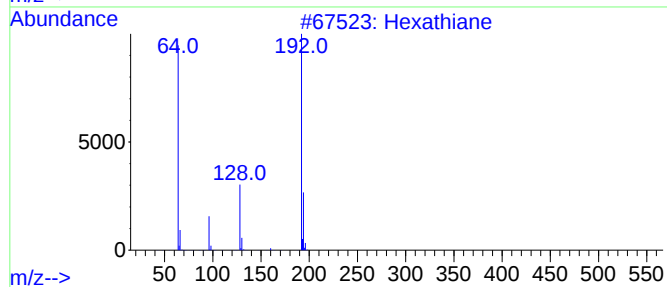
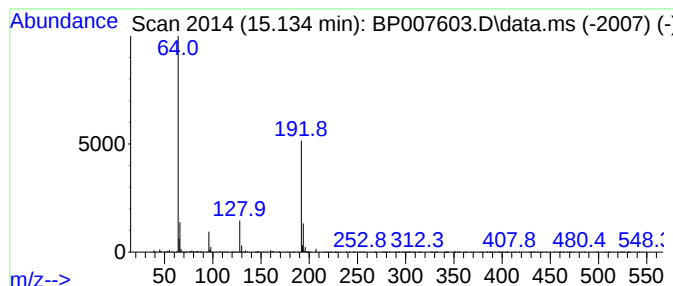
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 Hexathiane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.134	3.29 ng/ul	400114	Acenaphthene-d10	14.581

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			Cyclic octaatomic sulfur	256	S8	010544-50-0	56
5			Cyclobutane, 1,1'-(1,1,2,2-tetra...	386	C10H6Cl2F10	035208-02-7	9



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102121\
Data File : BP007603.D
Acq On : 22 Oct 2021 23:10
Operator : CG/JU
Sample : M4281-05DL 5X
Misc :
ALS Vial : 63 Sample Multiplier: 1

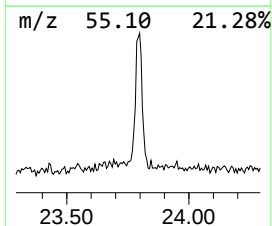
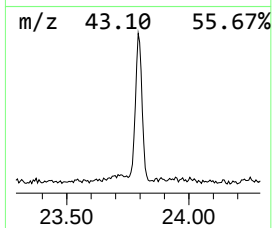
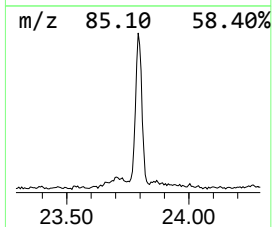
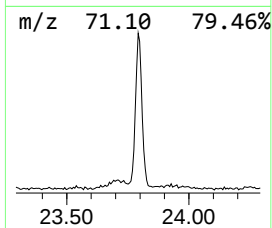
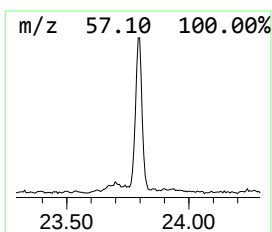
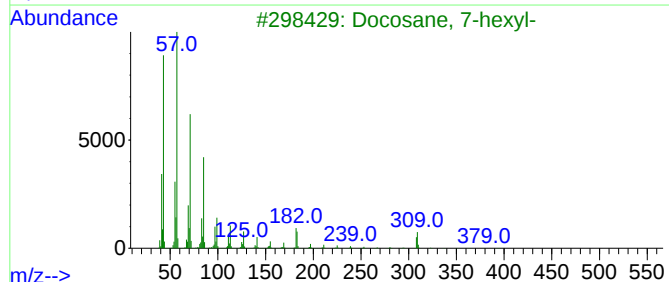
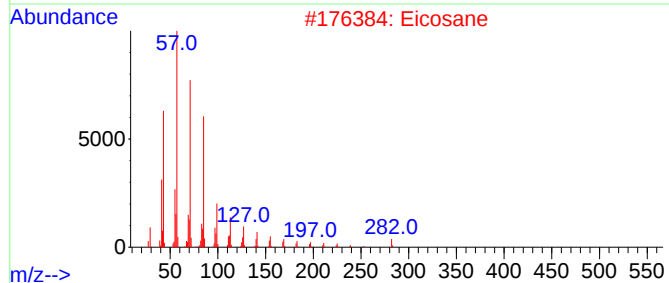
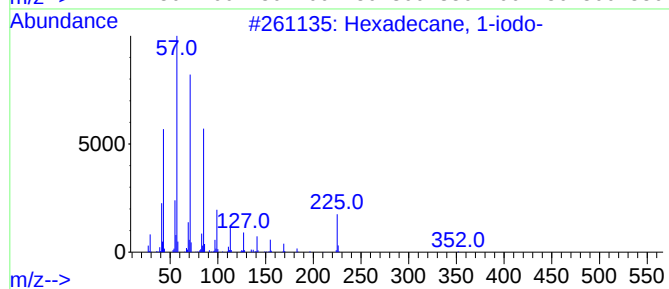
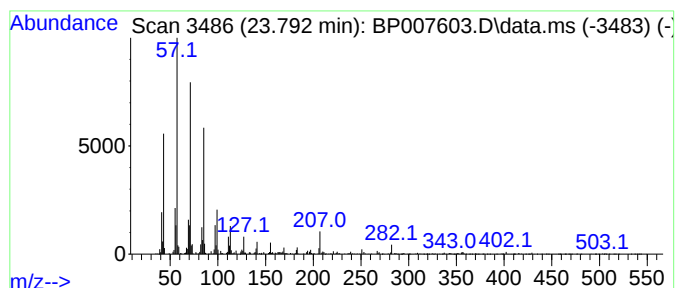
Instrument :
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ClientSampleId :
BFY61DL

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 11 Hexadecane, 1-iodo- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.	
23.792	2.02 ng/ul	360310	Perylene-d12	23.863	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane, 1-iodo-	352	C16H33I	000544-77-4	96
2	Eicosane	282	C20H42	000112-95-8	93
3	Docosane, 7-hexyl-	394	C28H58	055373-86-9	93
4	11-Methylpentacosane	366	C26H54	015689-71-1	93
5	Hexacosane	366	C26H54	000630-01-3	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102121\
 Data File : BP007603.D
 Acq On : 22 Oct 2021 23:10
 Operator : CG/JU
 Sample : M4281-05DL 5X
 Misc :
 ALS Vial : 63 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BFY61DL

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

TIC Library : C:\DATABASE\NIST0.L
 TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butanoic acid	4.052	10.2	ng/ul	595318	1	7.946	1163320	20.0
Butanoic acid, ...	4.817	4.4	ng/ul	254426	1	7.946	1163320	20.0
Butanoic acid, ...	5.017	16.3	ng/ul	946336	1	7.946	1163320	20.0
Pentanoic acid	5.517	31.8	ng/ul	1848490	1	7.946	1163320	20.0
Pentanoic acid,...	6.405	17.8	ng/ul	1037680	1	7.946	1163320	20.0
Hexanoic acid	7.034	25.5	ng/ul	1481470	1	7.946	1163320	20.0
Heptanoic acid	8.563	10.9	ng/ul	634435	1	7.946	1163320	20.0
Octanoic acid	10.105	10.3	ng/ul	889458	2	10.752	1725210	20.0
Ethanol, 2-(2-b...	10.534	3.0	ng/ul	262651	2	10.752	1725210	20.0
Hexathiane	15.134	3.3	ng/ul	400114	3	14.581	2435960	20.0
Hexadecane, 1-i...	23.792	2.0	ng/ul	360310	6	23.863	3564790	20.0