

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121022\
 Data File : BP013063.D
 Acq On : 11 Dec 2022 04:57
 Operator : CG/JU
 Sample : N5927-06
 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 EXZ92

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-BP120722.M
 Title : SVOA CALIBRATION

Signal : TIC: BP013063.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	24	28	39	rVB	85873	135643	0.54%	0.054%
2	3.781	98	101	118	rBV	730566	1210544	4.79%	0.480%
3	4.011	136	140	149	rVB	69134	110637	0.44%	0.044%
4	4.111	153	157	164	rVB	50552	77247	0.31%	0.031%
5	5.387	368	374	380	rVB	33387	50473	0.20%	0.020%
6	6.034	477	484	495	rVB	24942	50646	0.20%	0.020%
7	7.122	660	669	685	rVB	760662	1254386	4.96%	0.497%
8	7.293	691	698	707	rBV	2835420	4367560	17.27%	1.732%
9	7.499	726	733	744	rBV	2664357	4263587	16.86%	1.690%
10	7.969	806	813	818	rBV	2665953	4260597	16.85%	1.689%
11	8.028	818	823	835	rVB	1241788	1821815	7.20%	0.722%
12	8.675	925	933	941	rBV	2325251	3625141	14.34%	1.437%
13	8.799	949	954	961	rVB	109739	173558	0.69%	0.069%
14	9.140	1004	1012	1022	rBV	3275480	5438704	21.51%	2.156%
15	9.540	1075	1080	1090	rVB4	40148	95166	0.38%	0.038%
16	9.863	1126	1135	1145	rBV	2730316	4458015	17.63%	1.767%
17	10.398	1218	1226	1237	rBV	4222580	6971883	27.57%	2.764%
18	10.557	1247	1253	1262	rBV	77778	140293	0.55%	0.056%
19	10.781	1284	1291	1305	rBV	3877220	6738963	26.65%	2.672%
20	10.916	1305	1314	1330	rBV	3167476	5507393	21.78%	2.183%
21	11.316	1374	1382	1387	rBV	1077476	1689703	6.68%	0.670%
22	11.375	1387	1392	1411	rVB	814348	1437888	5.69%	0.570%
23	11.563	1419	1424	1428	rBV4	40389	88372	0.35%	0.035%
24	11.651	1434	1439	1447	rVB	20732	39981	0.16%	0.016%
25	11.728	1447	1452	1461	rVB10	21046	51120	0.20%	0.020%
26	12.363	1555	1560	1567	rBV2	67589	114253	0.45%	0.045%
27	13.010	1665	1670	1679	rVB2	81394	149403	0.59%	0.059%
28	13.216	1701	1705	1712	rVB7	25297	44570	0.18%	0.018%
29	13.445	1737	1744	1748	rVV3	43080	88597	0.35%	0.035%
30	13.498	1748	1753	1760	rVB	56321	94241	0.37%	0.037%
31	13.569	1760	1765	1771	rBV5	25727	55229	0.22%	0.022%
32	13.928	1822	1826	1833	rVB8	20294	38861	0.15%	0.015%
33	14.016	1833	1841	1847	rBV	9116294	12558996	49.66%	4.979%
34	14.134	1856	1861	1869	rVB	25584	48362	0.19%	0.019%
35	14.251	1875	1881	1884	rBV3	98436	184126	0.73%	0.073%
36	14.304	1884	1890	1898	rVV	9111923	13434031	53.13%	5.326%

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
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37	14.387	1901	1904	1910	rVV3	37586	68540	0.27%	0.027%
38	14.457	1913	1916	1920	rVV2	20787	37549	0.15%	0.015%
39	14.610	1935	1942	1952	rBV	6375870	9466067	37.43%	3.753%
40	14.704	1954	1958	1964	rVB8	26437	50067	0.20%	0.020%
41	14.804	1964	1975	1989	rBV	756831	1314884	5.20%	0.521%
42	14.922	1991	1995	1999	rBV	62736	83061	0.33%	0.033%
43	15.016	2006	2011	2017	rVB7	57169	93849	0.37%	0.037%
44	15.239	2044	2049	2055	rVB	66990	96715	0.38%	0.038%
45	15.345	2061	2067	2073	rBV	4351249	5718083	22.61%	2.267%
46	15.398	2073	2076	2081	rVB2	69221	94122	0.37%	0.037%
47	15.604	2103	2111	2124	rBV	12686011	18785556	74.29%	7.448%
48	15.716	2124	2130	2143	rVV	4829111	6507858	25.74%	2.580%
49	15.839	2145	2151	2157	rVB4	81660	168538	0.67%	0.067%
50	15.957	2168	2171	2175	rVB3	36123	45911	0.18%	0.018%
51	16.004	2175	2179	2183	rVB2	53508	73310	0.29%	0.029%
52	16.051	2184	2187	2193	rVB4	33178	46548	0.18%	0.018%
53	16.110	2193	2197	2200	rBV	54019	71914	0.28%	0.029%
54	16.145	2200	2203	2213	rVB8	34383	74163	0.29%	0.029%
55	16.239	2215	2219	2222	rVV5	53455	77974	0.31%	0.031%
56	16.275	2222	2225	2228	rVV2	84373	120491	0.48%	0.048%
57	16.310	2228	2231	2239	rVV3	69686	155663	0.62%	0.062%
58	16.381	2239	2243	2248	rVB4	44573	71023	0.28%	0.028%
59	16.492	2258	2262	2266	rBV2	60742	81315	0.32%	0.032%
60	16.657	2286	2290	2297	rVB8	34818	72709	0.29%	0.029%
61	16.739	2302	2304	2309	rVB4	35859	43480	0.17%	0.017%
62	16.833	2316	2320	2325	rVB6	61786	93784	0.37%	0.037%
63	17.104	2361	2366	2370	rBV3	66500	120367	0.48%	0.048%
64	17.157	2370	2375	2382	rVB2	214806	330505	1.31%	0.131%
65	17.233	2383	2388	2392	rBV	302830	459712	1.82%	0.182%
66	17.269	2392	2394	2400	rVB2	157672	201978	0.80%	0.080%
67	17.363	2404	2410	2417	rBV	8292283	11659590	46.11%	4.623%
68	17.463	2421	2427	2437	rVB	11908945	16858029	66.67%	6.684%
69	17.645	2453	2458	2468	rVV	160033	297017	1.17%	0.118%
70	17.739	2471	2474	2484	rVB6	64066	127531	0.50%	0.051%
71	18.016	2518	2521	2526	rVB	255128	325231	1.29%	0.129%
72	18.133	2537	2541	2547	rVB3	83220	141649	0.56%	0.056%
73	18.233	2553	2558	2562	rBV	978726	1250680	4.95%	0.496%
74	18.275	2562	2565	2567	rVV	250625	330867	1.31%	0.131%
75	18.304	2567	2570	2580	rVB	362049	553678	2.19%	0.220%
76	18.563	2604	2614	2619	rBV3	240670	532664	2.11%	0.211%

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77	18.645	2624	2628	2632	rVB3	101409	116575	0.46%	0.046%
78	18.716	2635	2640	2646	rBV3	319116	461877	1.83%	0.183%
79	18.810	2653	2656	2660	rBV	77041	95885	0.38%	0.038%
80	18.969	2680	2683	2687	rVB	125348	145502	0.58%	0.058%
81	19.057	2694	2698	2702	rBV	168559	218725	0.86%	0.087%
82	19.122	2706	2709	2711	rBV2	74155	82788	0.33%	0.033%
83	19.269	2730	2734	2737	rBV2	243552	296442	1.17%	0.118%
84	19.333	2741	2745	2749	rBV	241389	369432	1.46%	0.146%
85	19.463	2763	2767	2776	rBV	908218	1249784	4.94%	0.495%
86	19.692	2800	2806	2817	rVB	18321228	25287418	100.00%	10.026%
87	20.686	2972	2975	2978	rVB	127214	144225	0.57%	0.057%
88	20.774	2987	2990	2994	rBV	143108	183741	0.73%	0.073%
89	20.821	2995	2998	3001	rVB2	124620	146949	0.58%	0.058%
90	21.139	3048	3052	3064	rVB	1172670	1624055	6.42%	0.644%
91	21.451	3099	3105	3112	rBV	9162503	12490564	49.39%	4.952%
92	21.586	3125	3128	3132	rVB	198580	236115	0.93%	0.094%
93	22.057	3206	3208	3217	rVB2	219413	294149	1.16%	0.117%
94	22.574	3292	3296	3304	rVB	411623	615196	2.43%	0.244%
95	23.145	3390	3393	3399	rVB	490701	770811	3.05%	0.306%
96	23.780	3492	3501	3514	rVB2	11603506	24918946	98.54%	9.880%
97	23.939	3520	3528	3544	rVB2	5684711	12119911	47.93%	4.805%
98	24.557	3629	3633	3642	rVB	666639	1359166	5.37%	0.539%
99	25.439	3779	3783	3792	rVB	606153	1334836	5.28%	0.529%
100	28.092	4224	4234	4251	rVB	2939858	10092129	39.91%	4.001%

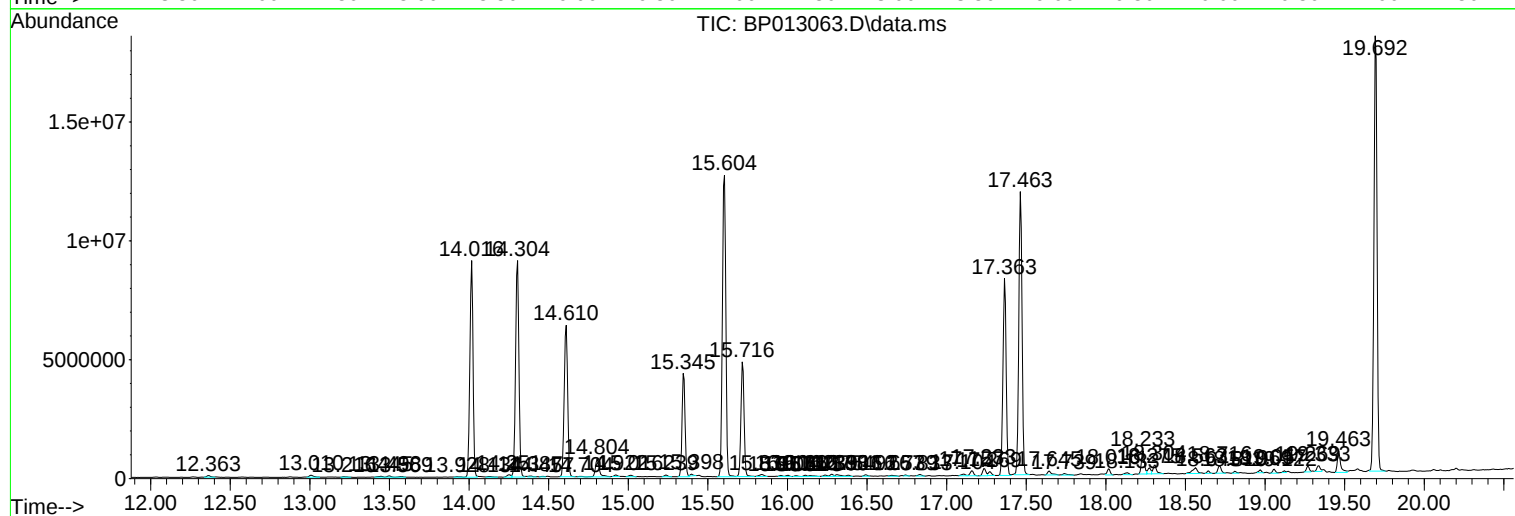
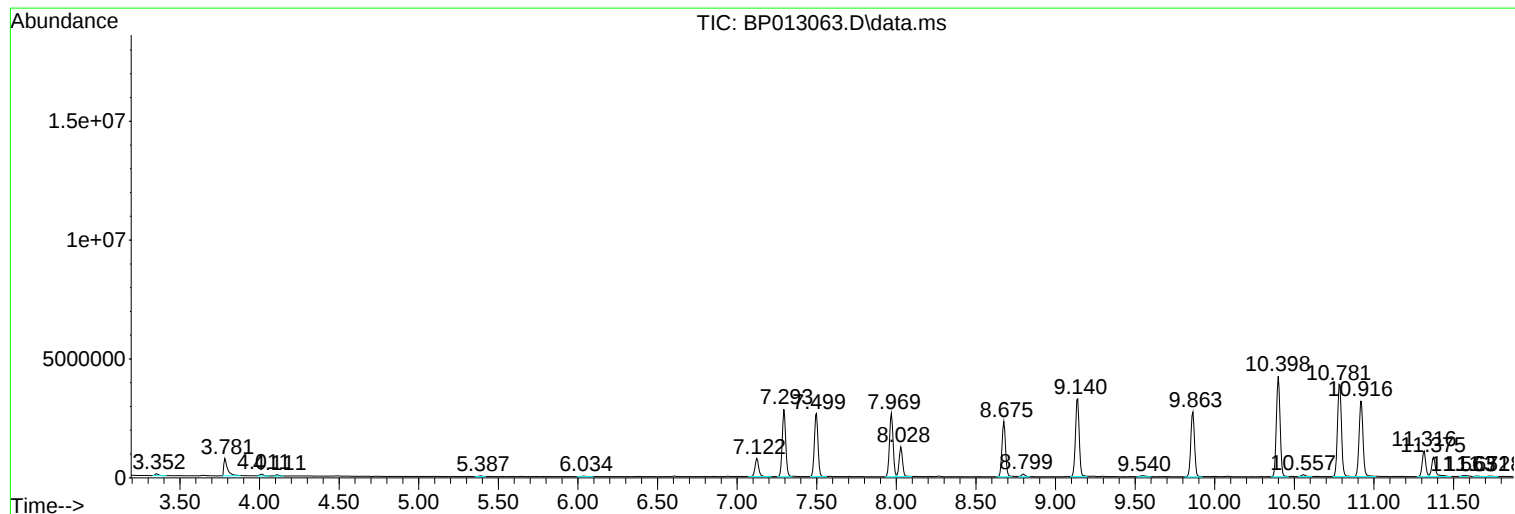
Sum of corrected areas: 252227877

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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



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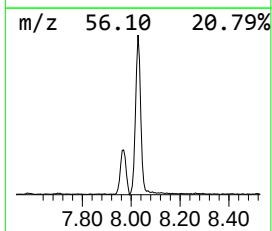
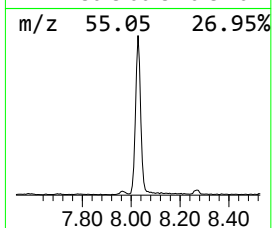
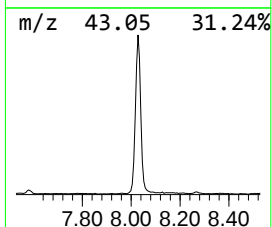
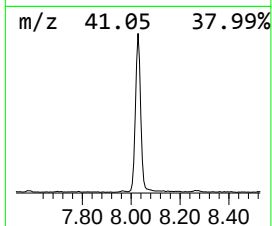
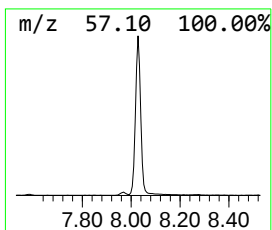
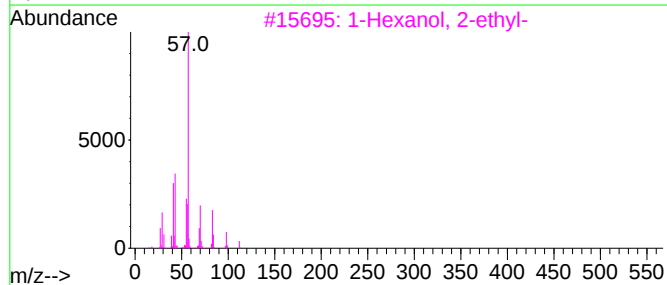
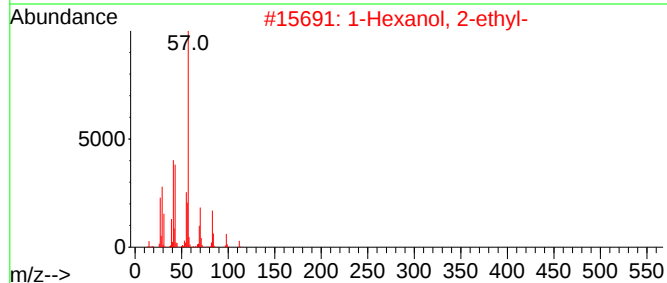
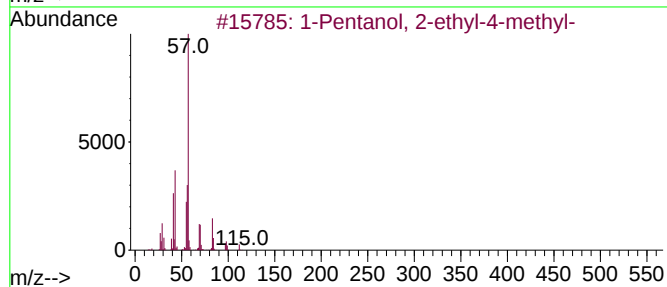
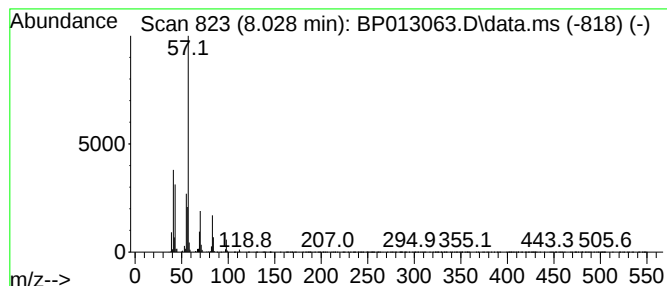
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 1-Pentanol, 2-ethyl-4-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.028	8.55 ng/ul	1821820	1,4-Dichlorobenzene-d4	7.969

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Pentanol, 2-ethyl-4-methyl-	130	C8H18O	000106-67-2	83		
2	1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	78		
3	1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	78		
4	1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	72		
5	1-Isobutoxy-2-ethylhexane	186	C12H26O	1000139-90-3	64		



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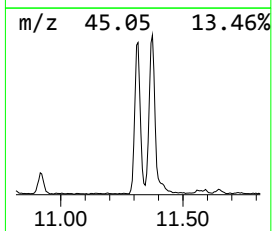
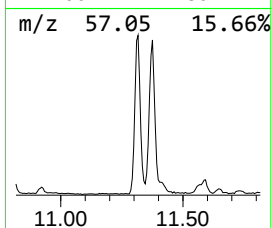
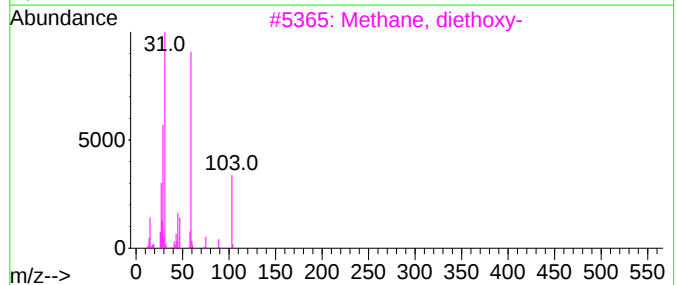
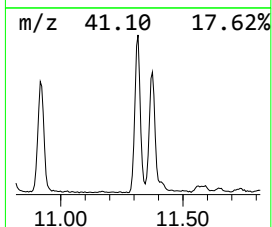
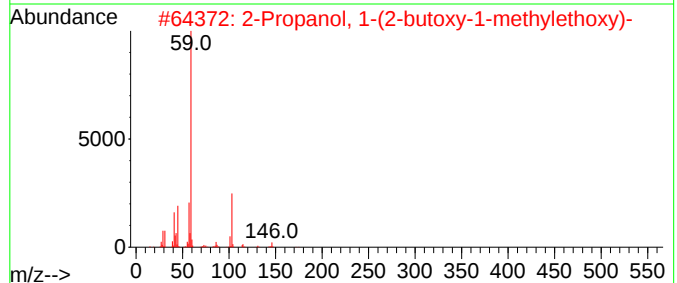
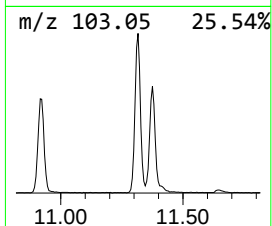
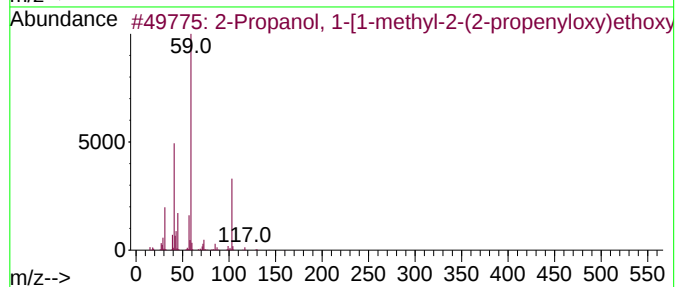
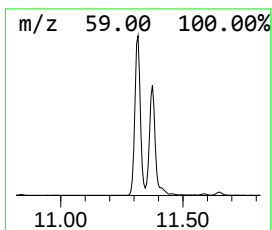
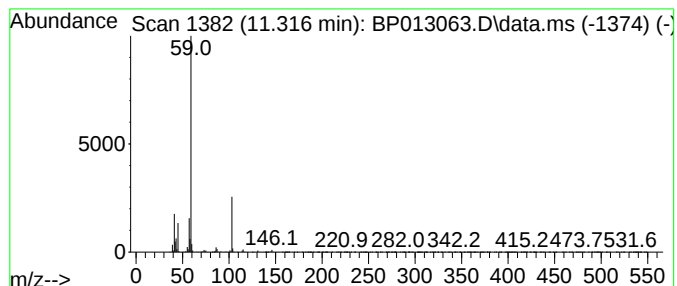
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TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 2-Propanol, 1-[1-methyl-2-(... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.316	5.01 ng/ul	1689700	Naphthalene-d8	10.781

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Propanol, 1-[1-methyl-2-(2-pro...	174	C9H18O3	055956-25-7	83		
2	2-Propanol, 1-(2-butoxy-1-methyl...	190	C10H22O3	029911-28-2	83		
3	Methane, diethoxy-	104	C5H12O2	000462-95-3	80		
4	1-Propanol, 2,2'-oxybis-	134	C6H14O3	000108-61-2	64		
5	1-Propanol, 2-(2-hydroxypropoxy)-	134	C6H14O3	000106-62-7	64		



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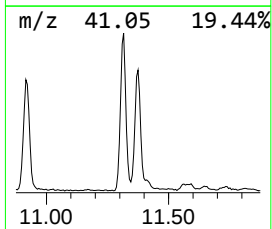
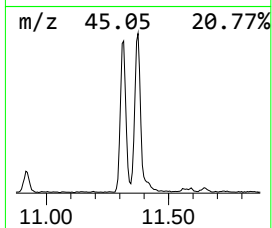
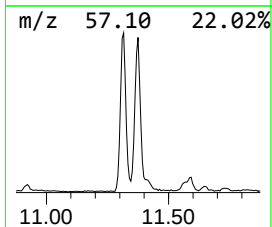
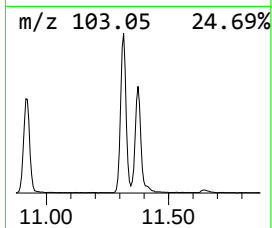
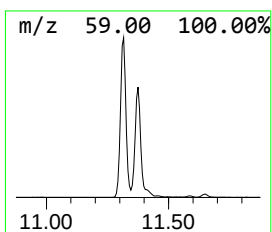
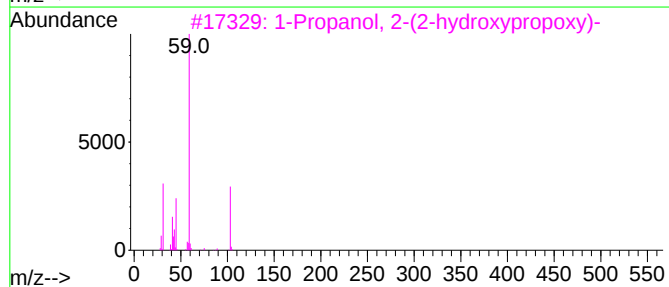
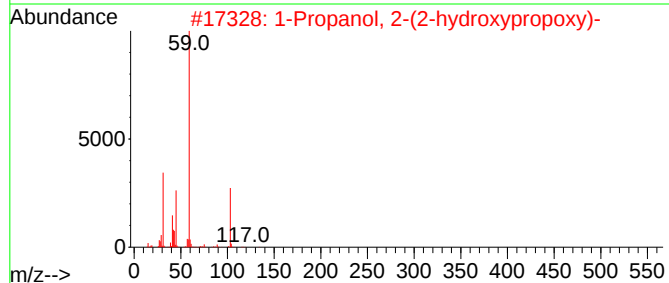
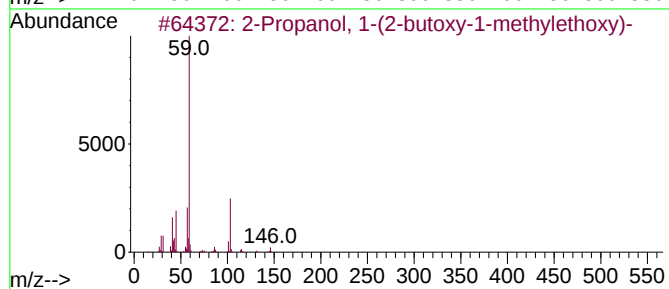
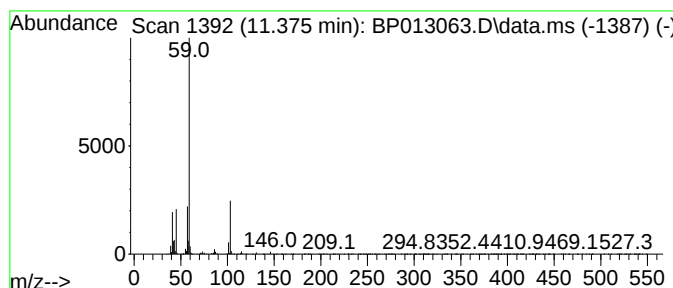
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 2-Propanol, 1-(2-butoxy-1-m... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.375	4.27 ng/ul	1437890	Naphthalene-d8	10.781

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Propanol, 1-(2-butoxy-1-methyl...	190	C10H22O3	029911-28-2	90		
2	1-Propanol, 2-(2-hydroxypropoxy)-	134	C6H14O3	000106-62-7	78		
3	1-Propanol, 2-(2-hydroxypropoxy)-	134	C6H14O3	000106-62-7	78		
4	2-Propanol, 1-[1-methyl-2-(2-pro...	174	C9H18O3	055956-25-7	64		
5	Dipropylene glycol (isomer 3)	134	C6H14O3	1000506-25-5	64		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121022\
Data File : BP013063.D
Acq On : 11 Dec 2022 04:57
Operator : CG/JU
Sample : N5927-06
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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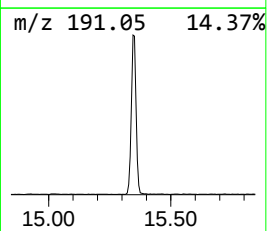
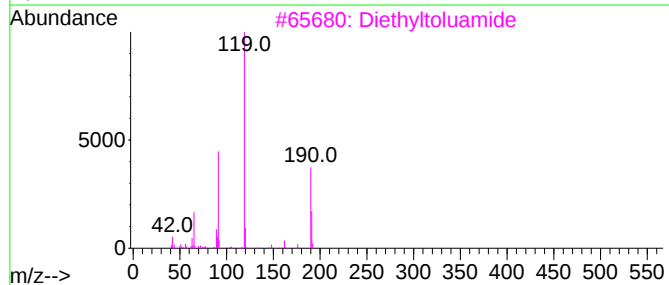
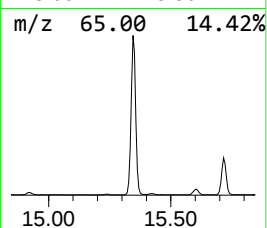
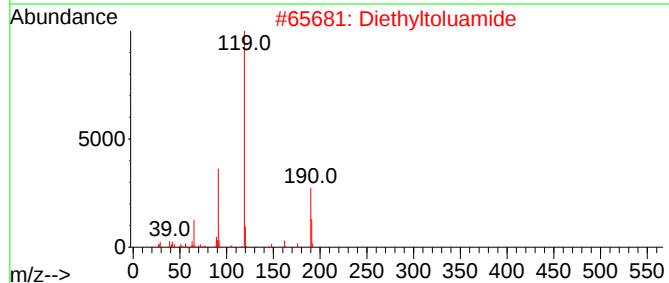
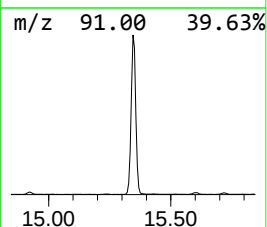
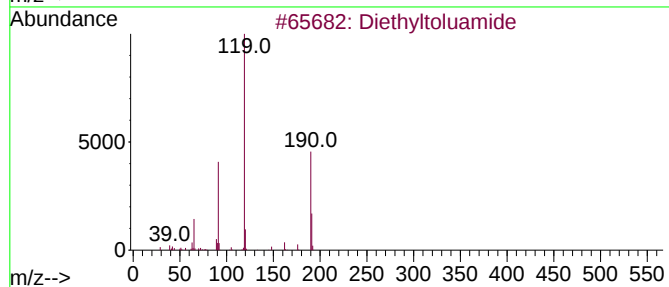
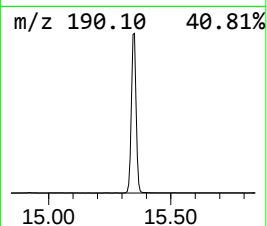
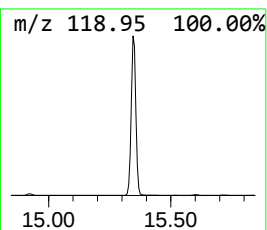
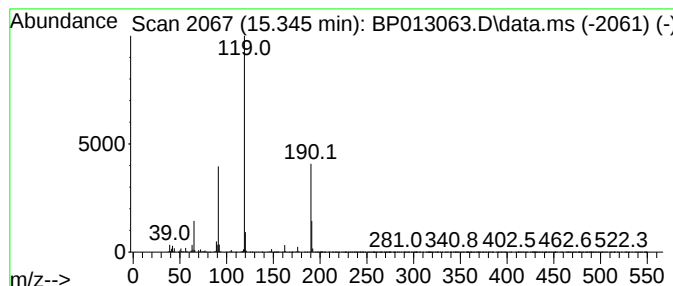
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Diethyltoluamide Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.345	12.08 ng/ul	5718080	Acenaphthene-d10	14.610

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Diethyltoluamide	191	C12H17NO	000134-62-3	97
2			Diethyltoluamide	191	C12H17NO	000134-62-3	95
3			Diethyltoluamide	191	C12H17NO	000134-62-3	94
4			Diethyltoluamide	191	C12H17NO	000134-62-3	93
5			Benzamide, N,N-diethyl-4-methyl-	191	C12H17NO	002728-05-4	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121022\
Data File : BP013063.D
Acq On : 11 Dec 2022 04:57
Operator : CG/JU
Sample : N5927-06
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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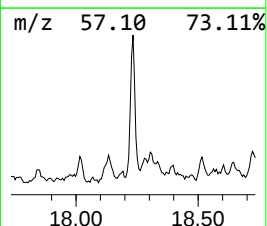
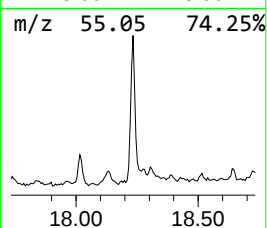
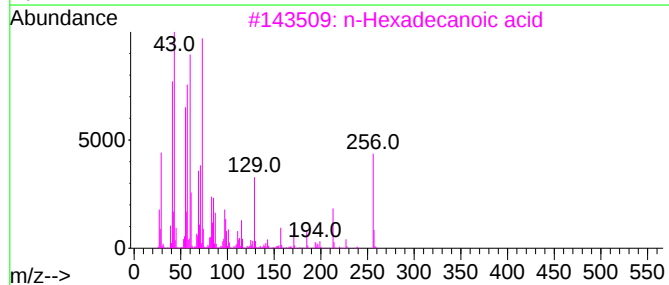
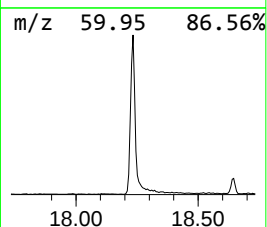
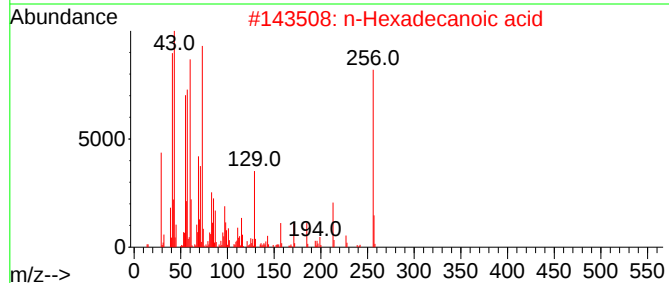
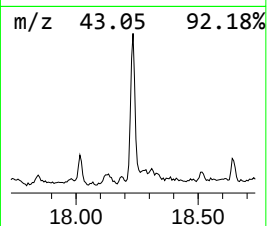
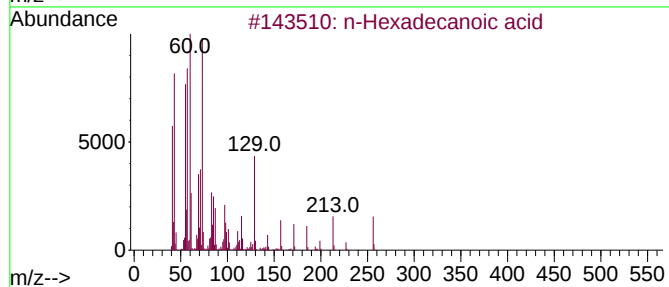
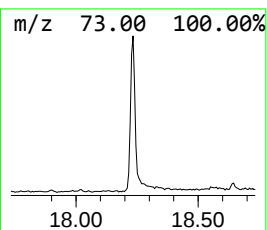
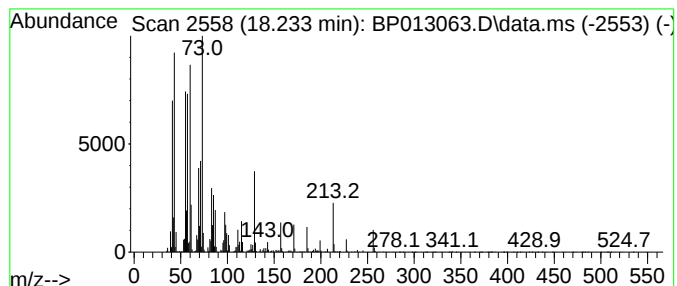
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 n-Hexadecanoic acid Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.233	2.15 ng/ul	1250680	Phenanthrene-d10	17.363

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	98
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
5			Octadecanoic acid	284	C18H36O2	000057-11-4	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121022\
Data File : BP013063.D
Acq On : 11 Dec 2022 04:57
Operator : CG/JU
Sample : N5927-06
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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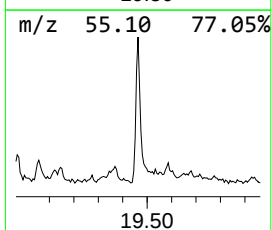
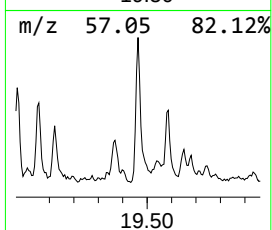
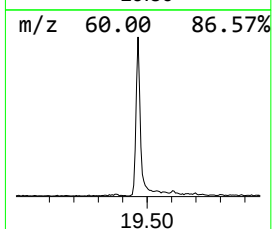
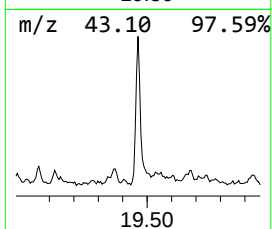
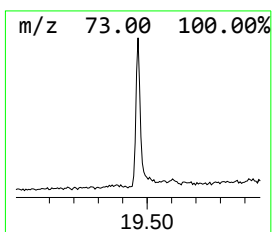
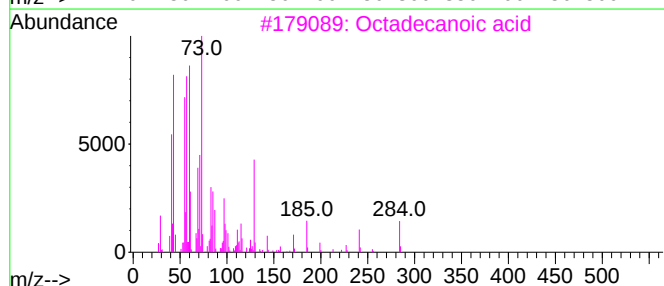
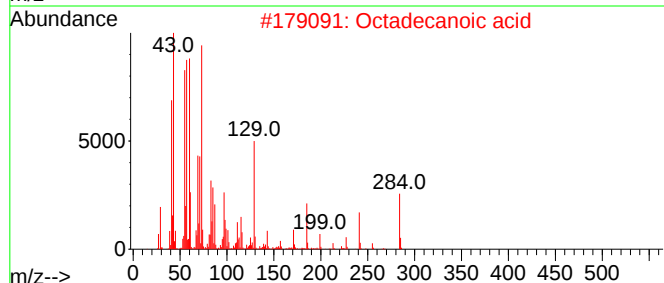
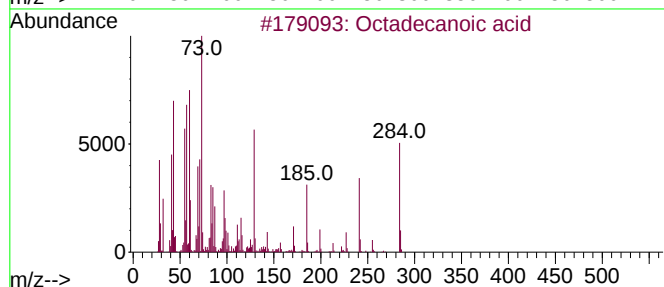
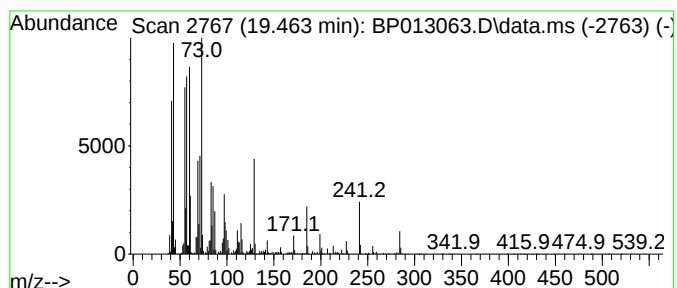
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Octadecanoic acid Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.463	2.00 ng/ul	1249780	Chrysene-d12	21.451

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	98
3			Octadecanoic acid	284	C18H36O2	000057-11-4	96
4			Octadecanoic acid	284	C18H36O2	000057-11-4	95
5			Octadecanoic acid	284	C18H36O2	000057-11-4	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121022\
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Acq On : 11 Dec 2022 04:57
Operator : CG/JU
Sample : N5927-06
Misc :
ALS Vial : 34 Sample Multiplier: 1

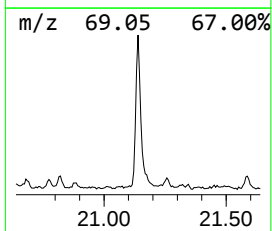
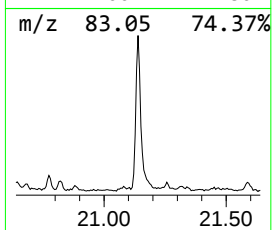
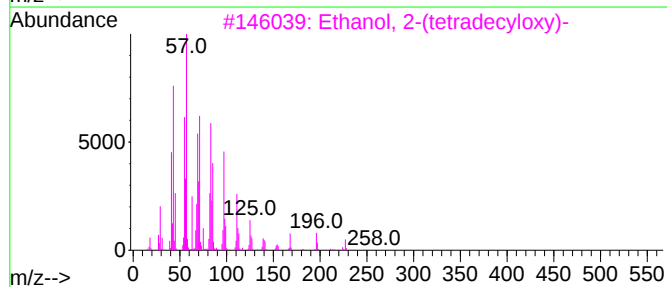
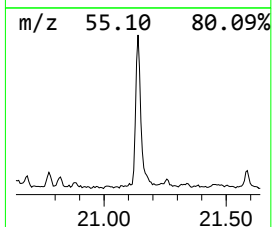
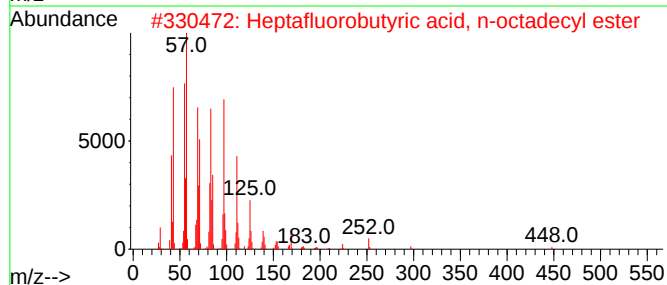
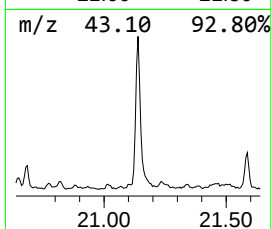
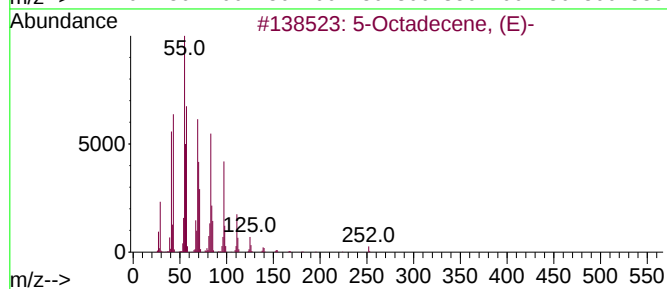
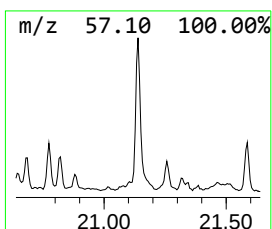
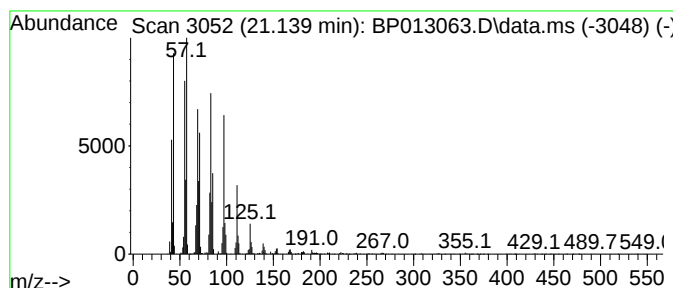
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
21.139	2.60 ng/ul	1624060	Chrysene-d12	21.451	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	5-Octadecene, (E)-	252	C18H36	007206-21-5	95
2	Heptafluorobutyric acid, n-octad...	466	C22H37F7O2	000400-57-7	94
3	Ethanol, 2-(tetradecyloxy)-	258	C16H34O2	002136-70-1	93
4	Cyclohexadecane	224	C16H32	000295-65-8	93
5	Pentadecafluorooctanoic acid, oc...	666	C26H37F15O2	1000406-04-8	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121022\
Data File : BP013063.D
Acq On : 11 Dec 2022 04:57
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Sample : N5927-06
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Instrument :
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ClientSampleId :
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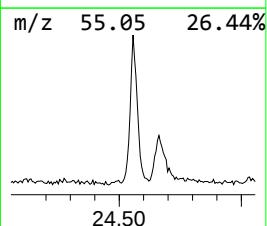
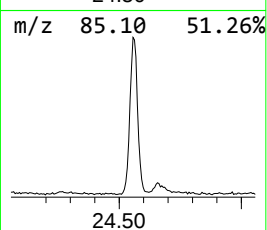
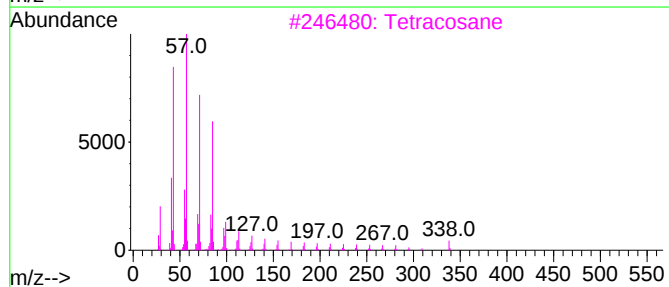
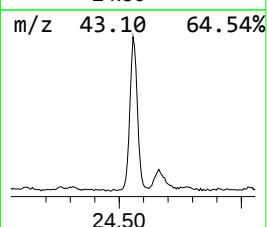
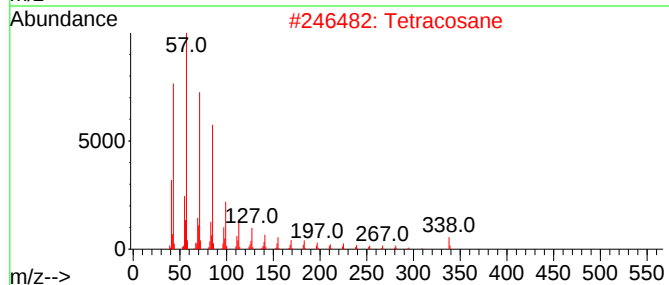
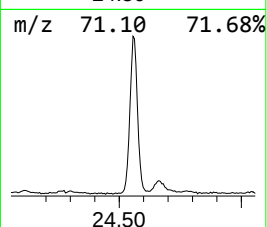
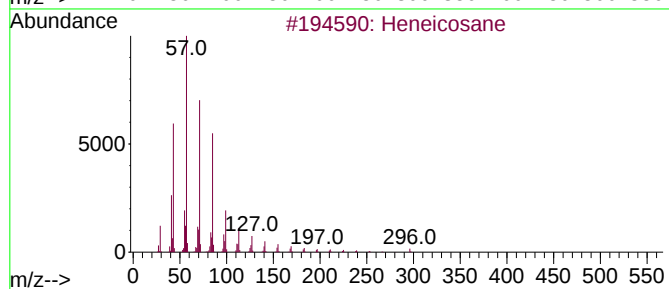
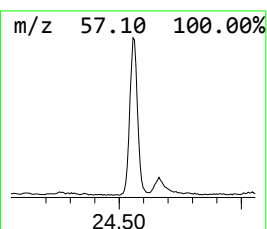
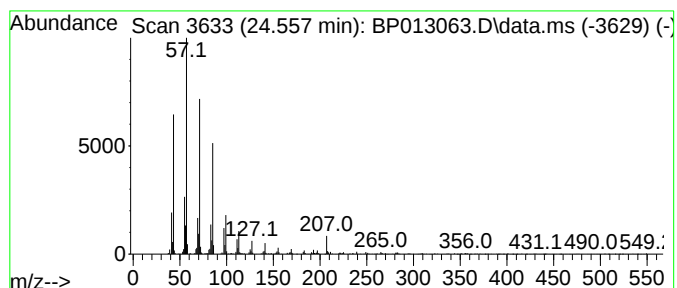
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.557	2.24 ng/ul	1359170	Perylene-d12	23.939

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heneicosane	296	C21H44	000629-94-7	97
2			Tetracosane	338	C24H50	000646-31-1	95
3			Tetracosane	338	C24H50	000646-31-1	93
4			2-Methyltetracosane	352	C25H52	001560-78-7	93
5			Eicosane	282	C20H42	000112-95-8	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121022\
Data File : BP013063.D
Acq On : 11 Dec 2022 04:57
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Misc :
ALS Vial : 34 Sample Multiplier: 1

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ClientSampleId :
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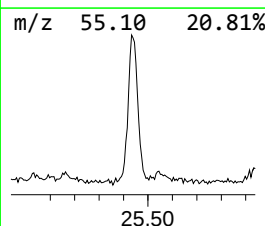
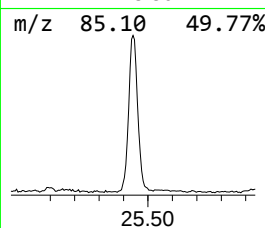
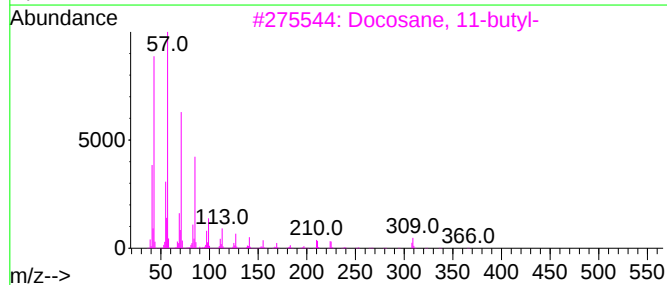
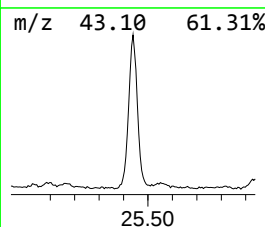
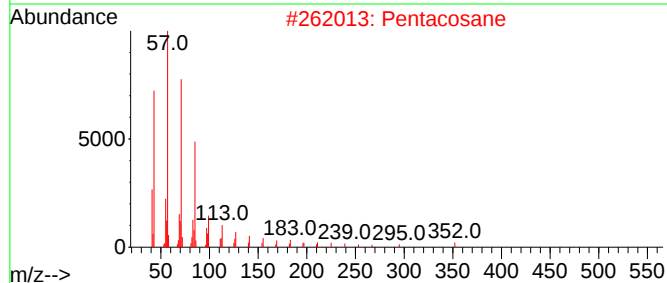
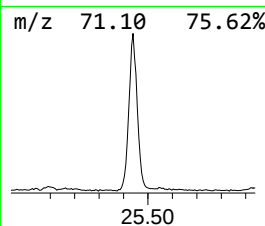
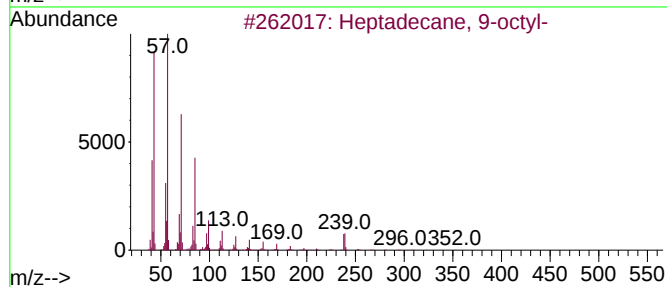
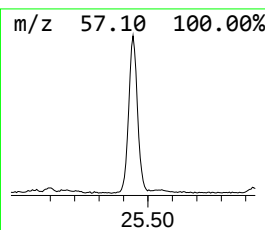
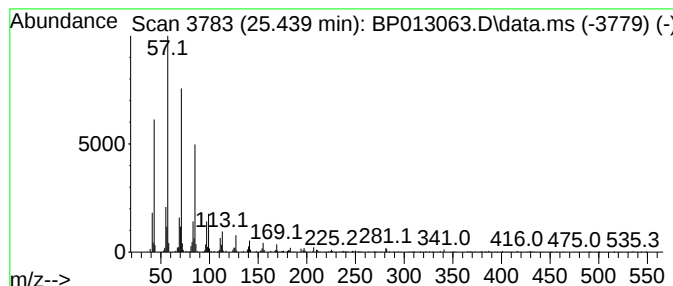
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.439	2.20 ng/ul	1334840	Perylene-d12	23.939

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	95
2		Pentacosane	352	C25H52	000629-99-2	94
3		Docosane, 11-butyl-	366	C26H54	013475-76-8	94
4		Octacosane, 2-methyl-	408	C29H60	001560-98-1	94
5		Eicosane	282	C20H42	000112-95-8	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121022\
Data File : BP013063.D
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Sample : N5927-06
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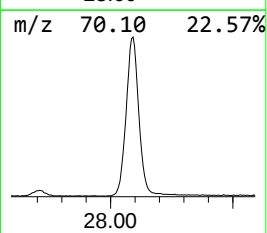
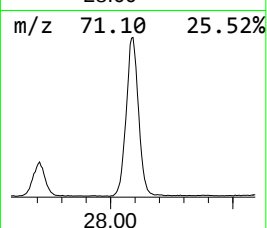
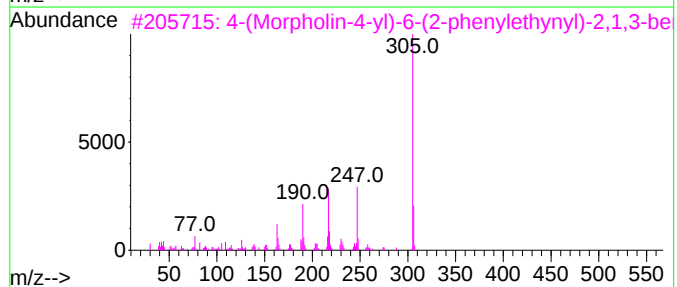
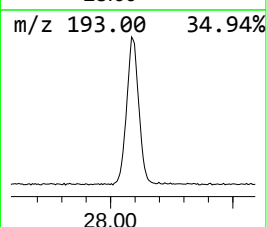
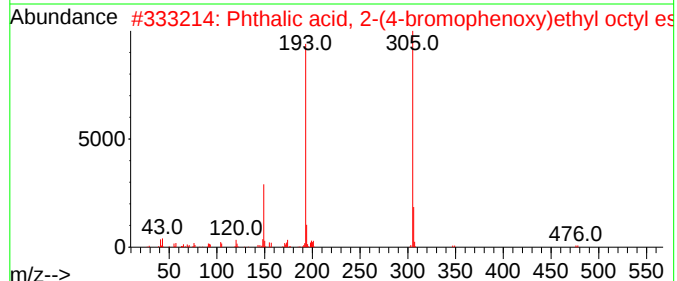
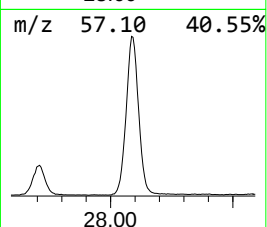
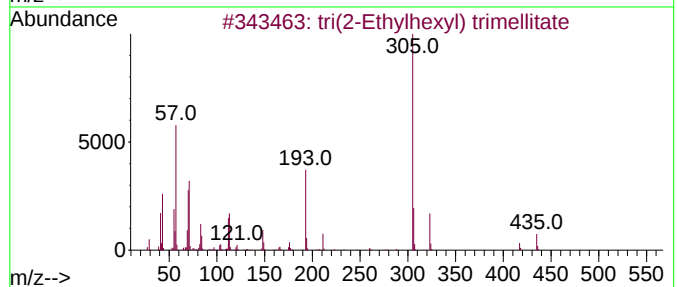
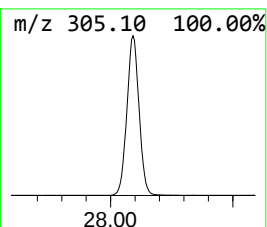
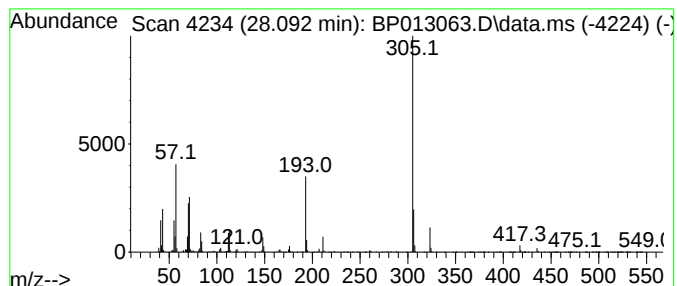
Instrument :
BNA_P
ClientSampleId :
EXZ92

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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

Peak Number 10 tri(2-Ethylhexyl) trimellitate Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.	
28.092	16.65 ng/ul	10092100	Perylene-d12	23.939	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	tri(2-Ethylhexyl) trimellitate	546	C33H54O6	003319-31-1	70
2	Phthalic acid, 2-(4-bromophenoxy...	476	C24H29BrO5	1010382-89-7	50
3	4-(Morpholin-4-yl)-6-(2-phenylet...	305	C18H15N3O2	1000387-57-0	23
4	4-(Diphenylphosphino)-N,N-dimeth...	305	C20H20NP	000739-58-2	23
5	(E)-N-(5-Methyl-1H-pyrazol-3-yl)...	375	C21H25N7	934353-76-1	23



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121022\
Data File : BP013063.D
Acq On : 11 Dec 2022 04:57
Operator : CG/JU
Sample : N5927-06
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EXZ92

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.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
1-Pentanol, 2-e...	8.028	8.6	ng/ul	1821820	1	7.969	4260600	20.0
2-Propanol, 1-[...	11.316	5.0	ng/ul	1689700	2	10.781	6738960	20.0
2-Propanol, 1-(...	11.375	4.3	ng/ul	1437890	2	10.781	6738960	20.0
Diethyltoluamide	15.345	12.1	ng/ul	5718080	3	14.610	9466070	20.0
n-Hexadecanoic ...	18.233	2.1	ng/ul	1250680	4	17.363	11659600	20.0
Octadecanoic acid	19.463	2.0	ng/ul	1249780	5	21.451	12490600	20.0
(DEL) Alkane: S...	21.139	2.6	ng/ul	1624060	5	21.451	12490600	20.0
(DEL) Alkane: S...	24.557	2.2	ng/ul	1359170	6	23.939	12119900	20.0
(DEL) Alkane: S...	25.439	2.2	ng/ul	1334840	6	23.939	12119900	20.0
tri(2-Ethylhexy...	28.092	16.6	ng/ul	10092100	6	23.939	12119900	20.0