

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102121\
 Data File : BP007552.D
 Acq On : 21 Oct 2021 13:29
 Operator : CG/JU
 Sample : M4262-04
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BFY68

Integration Parameters: LSCINT.P

Integrator: RTE

Soothing : 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: BP007552.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.375	12	15	20	rVV	35835	42427	0.82%	0.081%
2	3.658	59	63	70	rBV4	12639	19587	0.38%	0.037%
3	3.728	72	75	80	rVB	19274	22205	0.43%	0.042%
4	3.799	83	87	100	rBV	293770	479006	9.31%	0.910%
5	4.040	123	128	134	rVB2	27759	43585	0.85%	0.083%
6	4.128	140	143	149	rVB	12726	16249	0.32%	0.031%
7	4.228	156	160	164	rVB5	12437	17524	0.34%	0.033%
8	4.593	220	222	228	rVB4	7501	9645	0.19%	0.018%
9	4.781	249	254	261	rBV2	11688	20204	0.39%	0.038%
10	4.922	274	278	285	rBV2	16339	28937	0.56%	0.055%
11	5.052	296	300	304	rVB3	5395	9737	0.19%	0.018%
12	5.375	352	355	365	rVB8	6248	11447	0.22%	0.022%
13	5.793	422	426	430	rBV2	6522	10330	0.20%	0.020%
14	5.840	430	434	438	rVB3	8619	12993	0.25%	0.025%
15	6.034	463	467	475	rVB3	9221	12266	0.24%	0.023%
16	6.328	510	517	522	rBV5	12371	25327	0.49%	0.048%
17	6.457	532	539	542	rBV9	5204	10185	0.20%	0.019%
18	7.110	643	650	659	rVB2	221616	385418	7.49%	0.732%
19	7.275	671	678	686	rBV	519919	860292	16.72%	1.634%
20	7.440	702	706	707	rVV4	11479	16336	0.32%	0.031%
21	7.481	707	713	721	rVV	546411	861151	16.73%	1.635%
22	7.551	721	725	736	rVB	45584	76368	1.48%	0.145%
23	7.746	750	758	766	rVB5	19301	37646	0.73%	0.071%
24	7.863	773	778	783	rBV4	10173	16804	0.33%	0.032%
25	7.951	786	793	800	rVV	687448	1077189	20.93%	2.046%
26	8.022	800	805	809	rVB2	11569	17252	0.34%	0.033%
27	8.198	831	835	843	rVB3	11353	23618	0.46%	0.045%
28	8.381	861	866	868	rBV5	6162	9905	0.19%	0.019%
29	8.475	878	882	886	rBV2	14938	25842	0.50%	0.049%
30	8.516	886	889	894	rVB3	10731	13326	0.26%	0.025%
31	8.657	900	913	918	rBV	565342	965099	18.75%	1.833%
32	8.710	918	922	930	rVB	371259	565322	10.98%	1.074%
33	9.104	982	989	998	rBV	645817	1051259	20.43%	1.996%
34	9.281	1015	1019	1025	rBV7	8398	14354	0.28%	0.027%
35	9.498	1053	1056	1061	rBV7	4201	10624	0.21%	0.020%
36	9.575	1061	1069	1076	rVB2	19622	40946	0.80%	0.078%

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Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0.1 % of largest Peak

Start Thrs: 0.2

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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
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37	9.716	1088	1093	1103	rVB9	8284	17207	0.33%	0.033%
38	9.834	1103	1113	1120	rBV	425175	716742	13.93%	1.361%
39	10.081	1144	1155	1164	rBV	74778	134187	2.61%	0.255%
40	10.175	1164	1171	1173	rVV7	5377	9710	0.19%	0.018%
41	10.216	1173	1178	1184	rVV	15204	36295	0.71%	0.069%
42	10.375	1197	1205	1213	rBV	743912	1241219	24.12%	2.357%
43	10.540	1226	1233	1250	rBV	641825	1050937	20.42%	1.996%
44	10.757	1261	1270	1278	rVB	902590	1555652	30.23%	2.954%
45	10.887	1284	1292	1301	rBV	783970	1334727	25.93%	2.535%
46	11.551	1400	1405	1412	rBV7	8335	16149	0.31%	0.031%
47	11.757	1436	1440	1445	rBV3	7523	13769	0.27%	0.026%
48	11.816	1447	1450	1458	rVB4	8410	18850	0.37%	0.036%
49	11.951	1467	1473	1477	rBV	110265	170389	3.31%	0.324%
50	11.998	1477	1481	1482	rVV	126090	175000	3.40%	0.332%
51	12.016	1482	1484	1488	rVV2	134961	190871	3.71%	0.362%
52	12.057	1488	1491	1502	rVB	112750	192908	3.75%	0.366%
53	12.198	1510	1515	1519	rBV2	43939	77823	1.51%	0.148%
54	12.239	1519	1522	1527	rVV3	56023	94425	1.83%	0.179%
55	12.286	1527	1530	1535	rVV2	22765	41755	0.81%	0.079%
56	12.345	1535	1540	1544	rVB	16625	30183	0.59%	0.057%
57	12.469	1557	1561	1566	rVB6	7257	12288	0.24%	0.023%
58	12.816	1614	1620	1624	rBV9	9214	18536	0.36%	0.035%
59	12.869	1624	1629	1636	rVV2	13433	23792	0.46%	0.045%
60	12.957	1640	1644	1648	rBV5	7419	10945	0.21%	0.021%
61	13.204	1683	1686	1691	rVB4	11990	15809	0.31%	0.030%
62	13.492	1731	1735	1739	rBV3	8983	14159	0.28%	0.027%
63	13.710	1766	1772	1775	rBV6	6711	9843	0.19%	0.019%
64	13.898	1798	1804	1809	rBV	13386	25930	0.50%	0.049%
65	13.992	1814	1820	1828	rVV	1605573	2219429	43.12%	4.215%
66	14.051	1828	1830	1835	rVV4	15261	21285	0.41%	0.040%
67	14.116	1837	1841	1846	rVB6	6065	11093	0.22%	0.021%
68	14.275	1853	1868	1877	rBV	1655070	2548323	49.51%	4.839%
69	14.469	1896	1901	1905	rBV4	13627	26193	0.51%	0.050%
70	14.586	1913	1921	1928	rBV	1464476	2117485	41.14%	4.021%
71	14.769	1939	1952	1957	rBV	128624	210274	4.09%	0.399%
72	14.816	1957	1960	1968	rVB10	10699	22282	0.43%	0.042%
73	15.004	1987	1992	1994	rBV5	23707	33410	0.65%	0.063%
74	15.033	1994	1997	2004	rVB3	32645	50467	0.98%	0.096%
75	15.133	2007	2014	2030	rBV	99085	237696	4.62%	0.451%
76	15.298	2036	2042	2048	rBV	41119	66492	1.29%	0.126%

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Integration Parameters: LSCINT.P

Integrator: RTE
 Smoothing : OFF Filtering: 5
 Sampling : 1 Min Area: 0.1 % of largest Peak
 Start Thrs: 0.2 Max Peaks: 100
 Stop Thrs : 0 Peak Location: TOP

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 Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP101421.M
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77	15.398	2053	2059	2064	rVV4	11835	20772	0.40%	0.039%
78	15.451	2065	2068	2072	rVB5	11363	15025	0.29%	0.029%
79	15.580	2083	2090	2099	rBV	2367599	3458024	67.19%	6.567%
80	15.686	2102	2108	2117	rBV	318561	472058	9.17%	0.896%
81	15.816	2125	2130	2135	rBV	30414	46922	0.91%	0.089%
82	16.027	2159	2166	2170	rBV5	17051	31892	0.62%	0.061%
83	16.745	2283	2288	2292	rBV5	27708	39202	0.76%	0.074%
84	17.092	2343	2347	2355	rBV3	38678	70960	1.38%	0.135%
85	17.339	2382	2389	2396	rBV	1767333	2593936	50.40%	4.926%
86	17.439	2399	2406	2415	rBV	2937731	4168850	81.00%	7.917%
87	17.527	2418	2421	2427	rVB6	20981	32121	0.62%	0.061%
88	17.874	2478	2480	2488	rVB5	10018	18197	0.35%	0.035%
89	18.021	2501	2505	2511	rVB	139195	171000	3.32%	0.325%
90	18.239	2536	2542	2550	rBV	1182324	1577939	30.66%	2.996%
91	19.151	2693	2697	2701	rBV3	88289	106709	2.07%	0.203%
92	19.251	2711	2714	2717	rBV	40213	47643	0.93%	0.090%
93	19.321	2722	2726	2728	rBV3	75935	101392	1.97%	0.193%
94	19.351	2728	2731	2734	rVV4	44366	53758	1.04%	0.102%
95	19.468	2746	2751	2766	rBV	600192	907079	17.62%	1.723%
96	19.668	2779	2785	2793	rBV	3755585	5072177	98.55%	9.632%
97	21.145	3032	3036	3041	rBV	441965	568133	11.04%	1.079%
98	21.421	3078	3083	3090	rBV	2283708	3075577	59.76%	5.840%
99	23.715	3466	3473	3483	rVB2	2557800	5146631	100.00%	9.773%
100	23.874	3493	3500	3513	rVB2	1510427	3188831	61.96%	6.056%

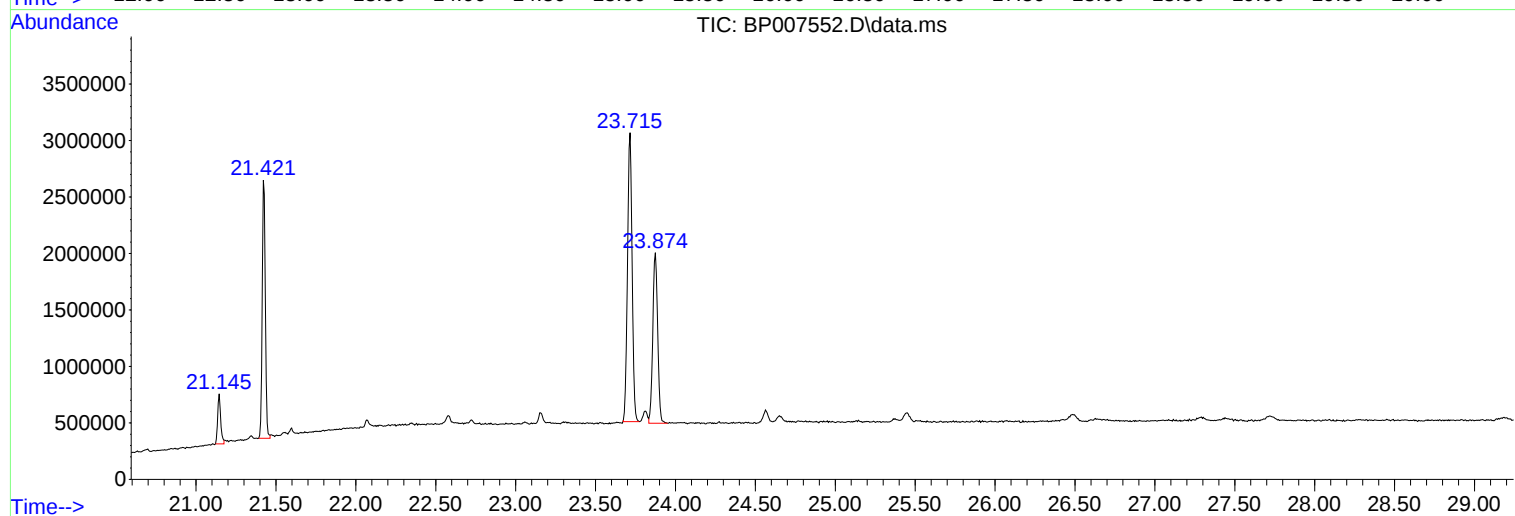
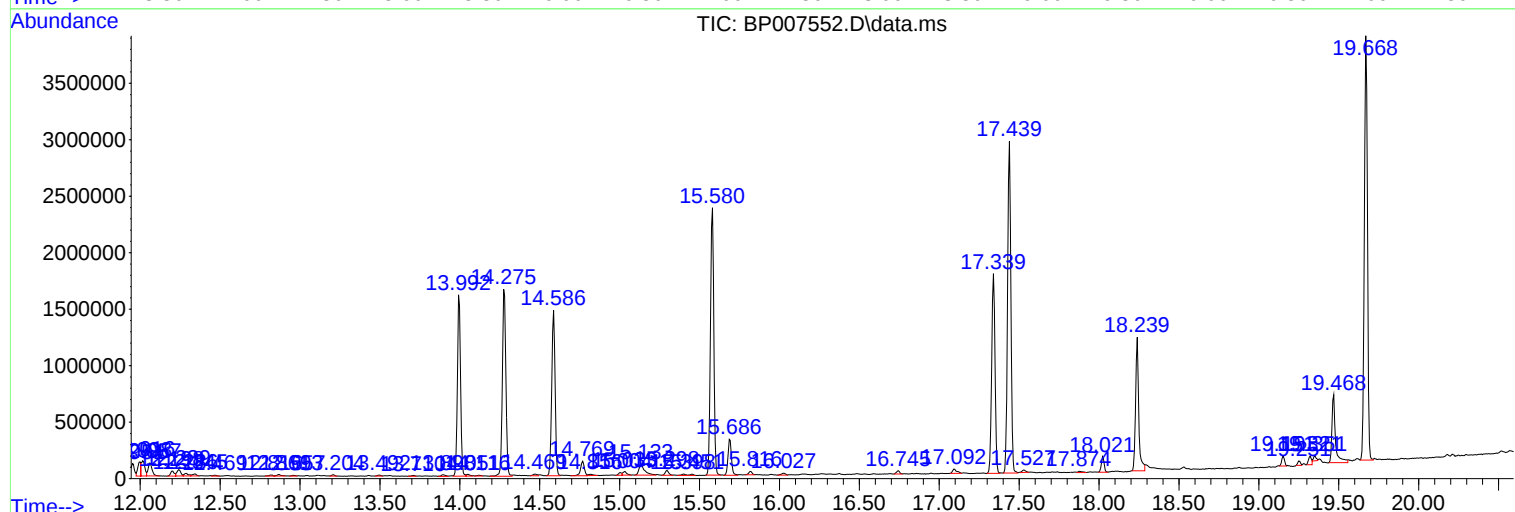
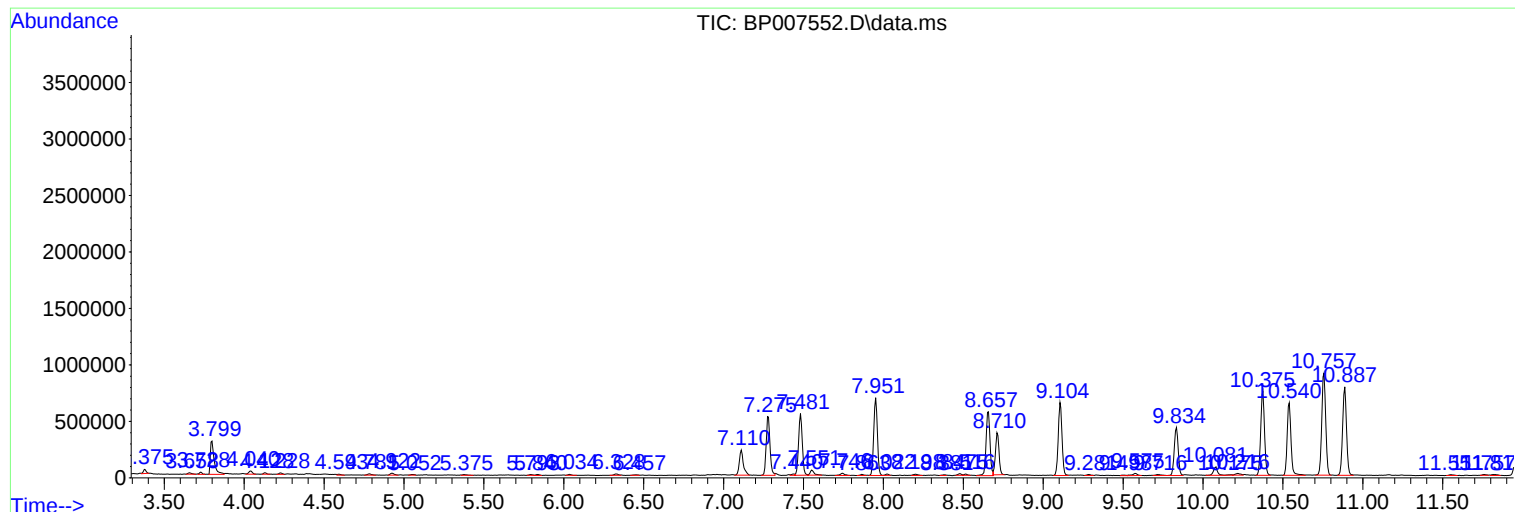
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Instrument :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP101421.M
Quant Title : SVOA CALIBRATION

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



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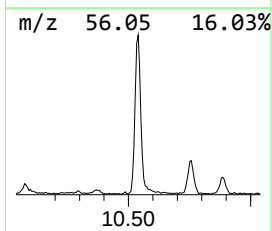
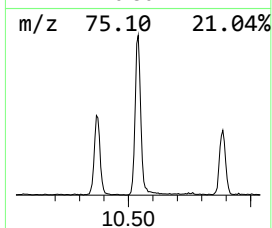
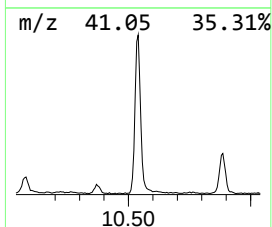
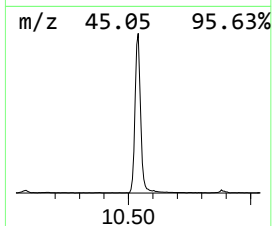
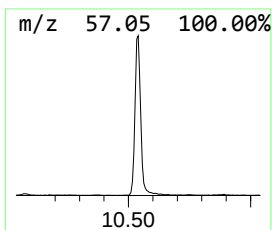
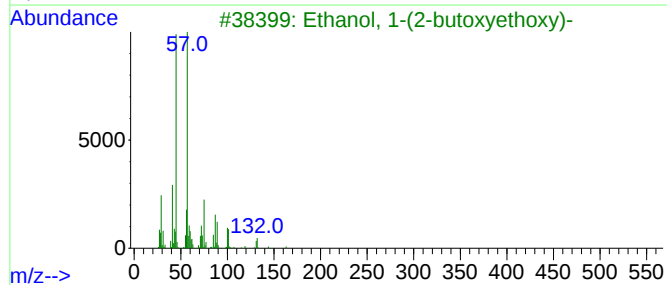
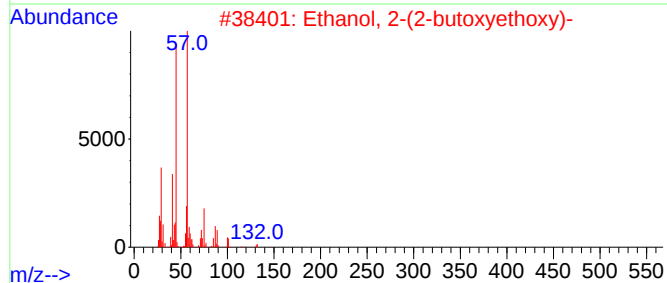
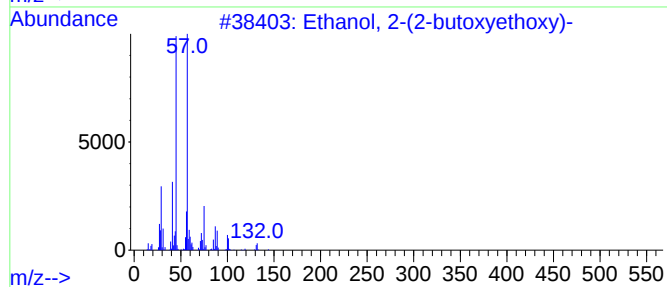
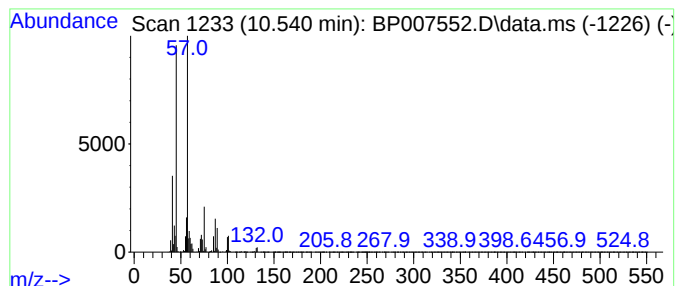
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP101421.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.540	13.51 ng/ul	1050940	Naphthalene-d8	10.757

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
2			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
3			Ethanol, 1-(2-butoxyethoxy)-	162	C8H18O3	054446-78-5	90
4			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
5			Ethanol, 2-[2-(2-methylpropoxy)e...	162	C8H18O3	018912-80-6	72



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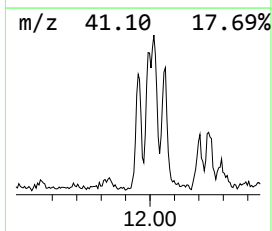
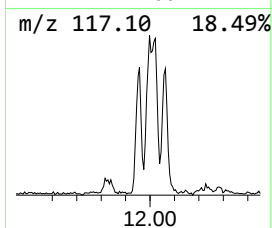
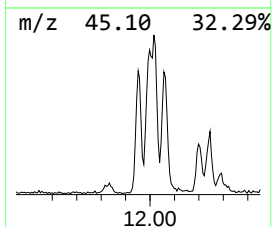
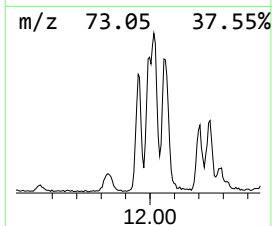
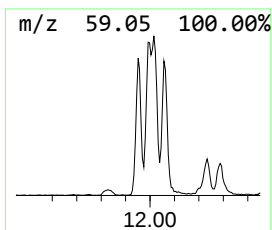
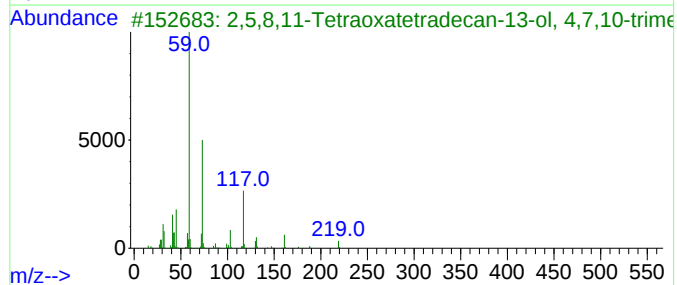
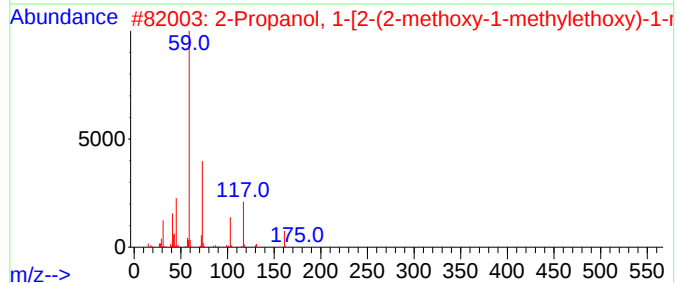
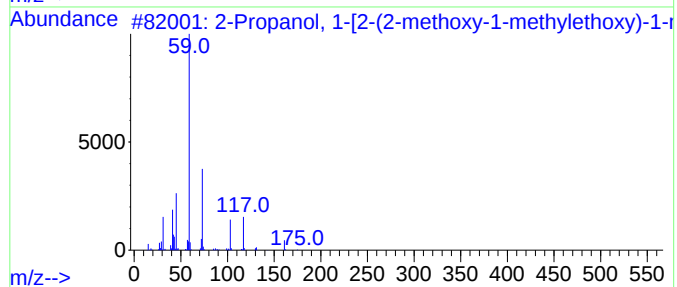
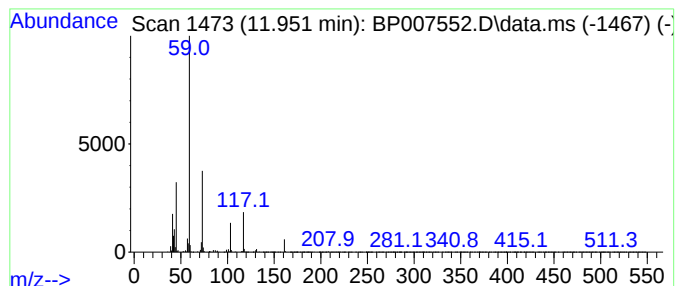
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Quant Title : SVOA CALIBRATION

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

Peak Number 2 2-Propanol, 1-[2-(2-methoxy... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.951	2.19 ng/ul	170389	Naphthalene-d8	10.757

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Propanol, 1-[2-(2-methoxy-1-me...	206	C10H22O4	020324-33-8	90		
2	2-Propanol, 1-[2-(2-methoxy-1-me...	206	C10H22O4	020324-33-8	86		
3	2,5,8,11-Tetraoxatetradecan-13-o...	264	C13H28O5	020324-34-9	72		
4	2-Propanol, 1-(2-methoxy-1-methy...	148	C7H16O3	020324-32-7	59		
5	1-Propanol, 2-(2-methoxypropoxy)-	148	C7H16O3	013588-28-8	59		



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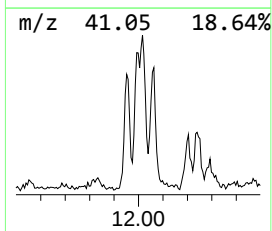
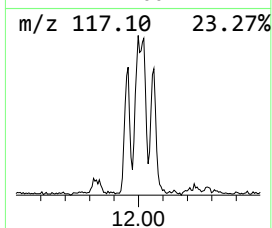
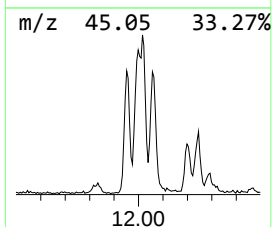
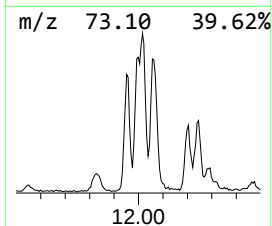
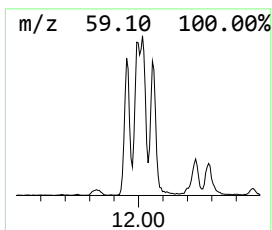
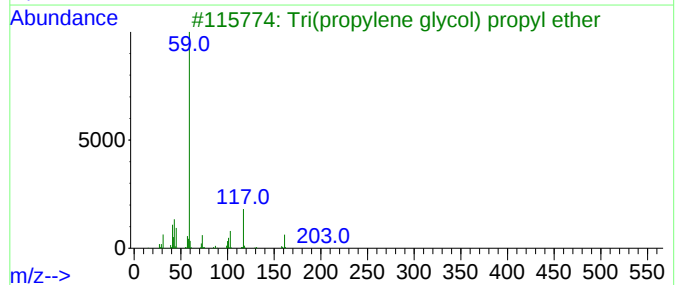
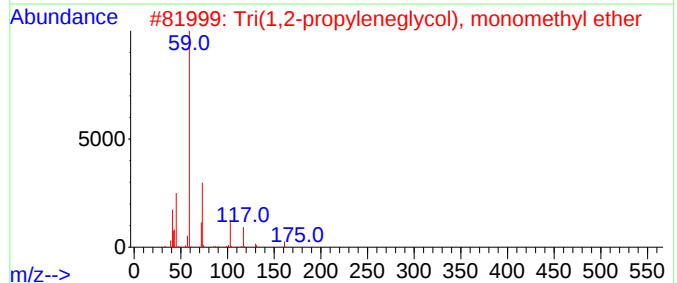
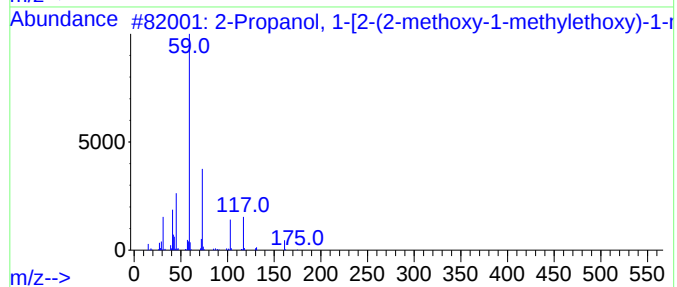
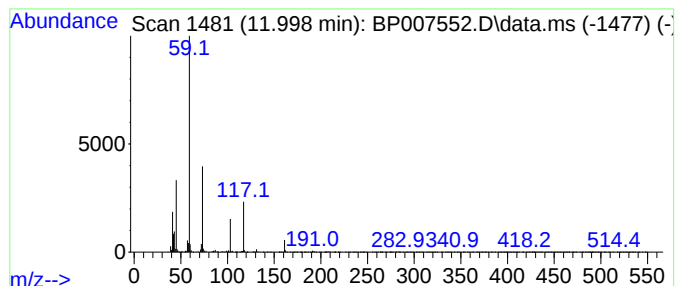
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP101421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 Tri(1,2-propyleneglycol), m... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.998	2.25 ng/ul	175000	Naphthalene-d8	10.757

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Propanol, 1-[2-(2-methoxy-1-me...	206	C10H22O4	020324-33-8	86		
2	Tri(1,2-propyleneglycol), monome...	206	C10H22O4	1000262-26-6	72		
3	Tri(propylene glycol) propyl ether	234	C12H26O4	096077-04-2	72		
4	2-Propanol, 1-(2-methoxy-1-methy...	148	C7H16O3	020324-32-7	64		
5	Methyl 2,5,8,11-tetraoxatridecan...	236	C10H20O6	1000366-78-2	56		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102121\
Data File : BP007552.D
Acq On : 21 Oct 2021 13:29
Operator : CG/JU
Sample : M4262-04
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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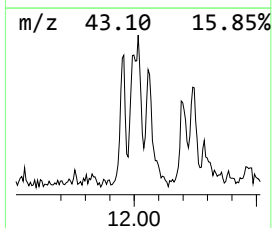
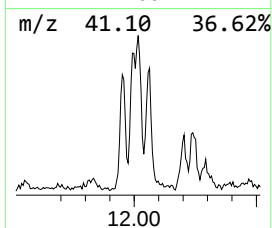
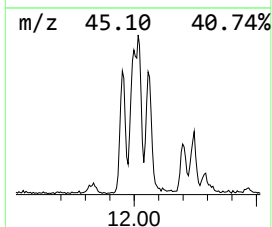
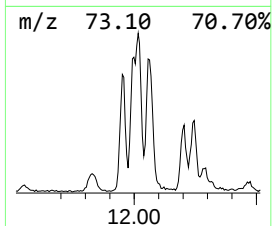
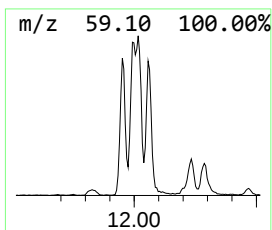
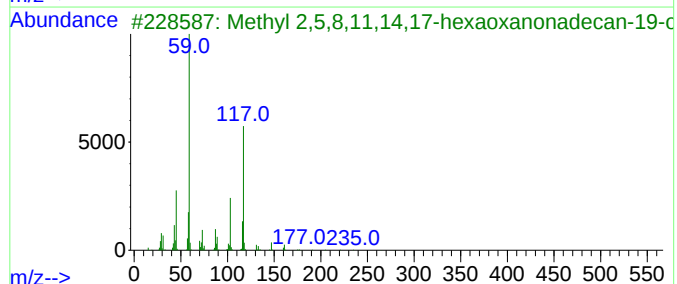
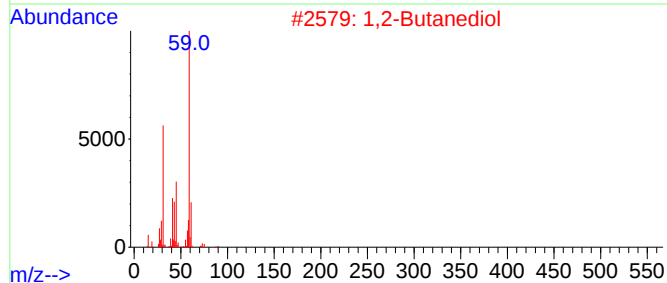
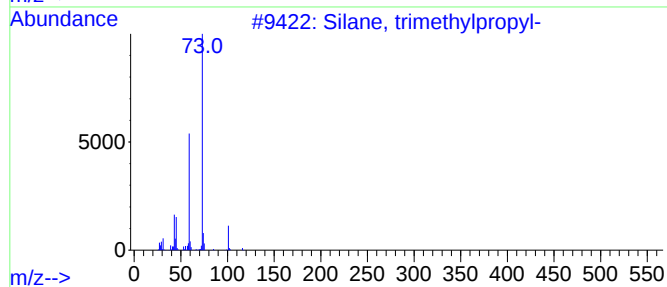
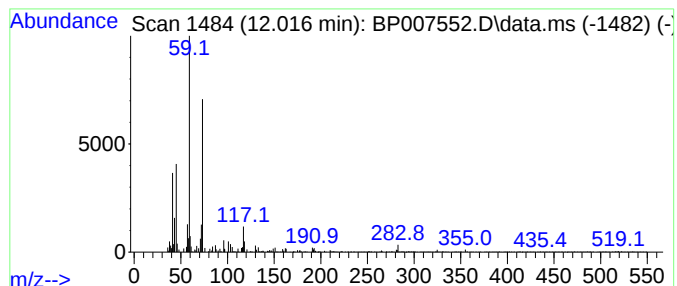
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Silane, trimethylpropyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.016	2.45 ng/ul	190871	Naphthalene-d8	10.757

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Silane, trimethylpropyl-	116	C6H16Si	003510-70-1	72
2			1,2-Butanediol	90	C4H10O2	000584-03-2	72
3			Methyl 2,5,8,11,14,17-hexaoxonon...	324	C14H28O8	1000366-81-4	56
4			Ethanedioic acid, dimethyl ester	118	C4H6O4	000553-90-2	56
5			Butanamide	87	C4H9NO	000541-35-5	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102121\
Data File : BP007552.D
Acq On : 21 Oct 2021 13:29
Operator : CG/JU
Sample : M4262-04
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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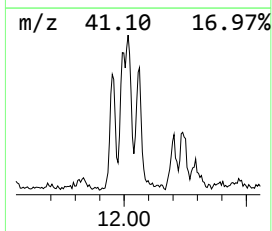
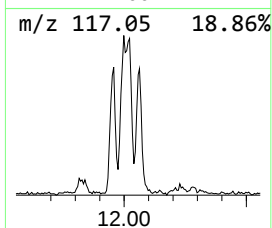
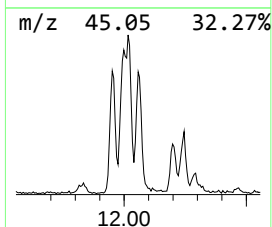
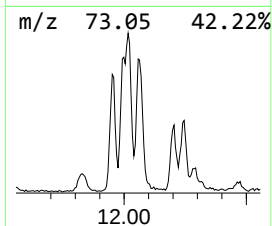
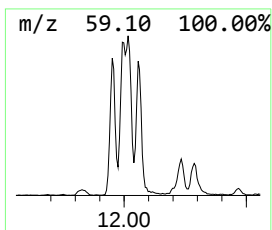
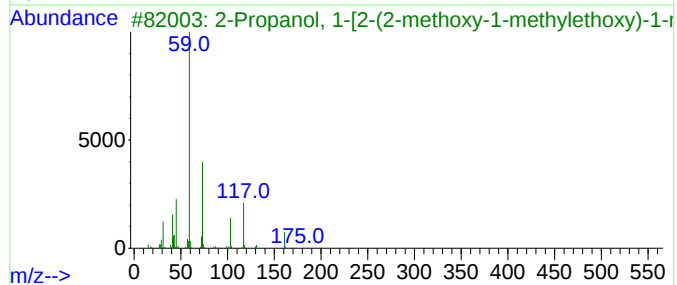
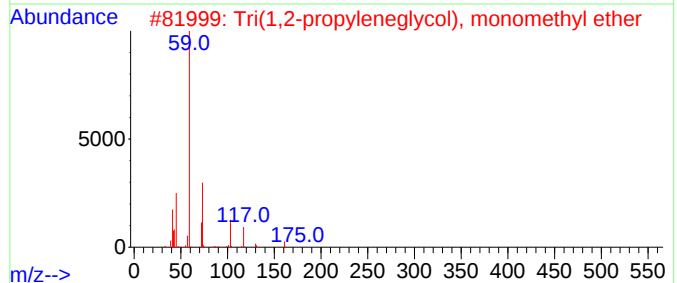
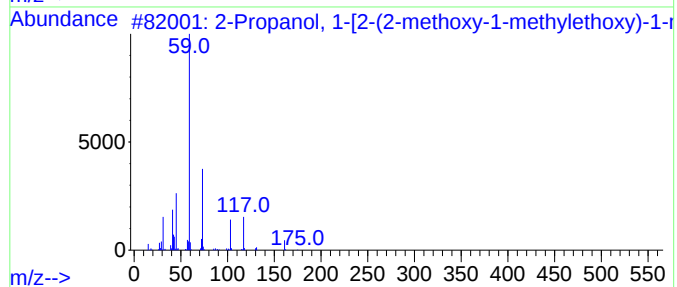
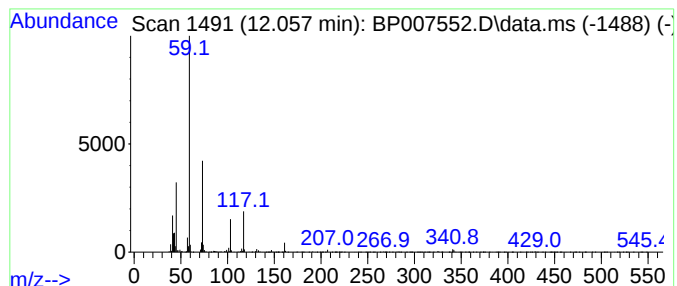
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Methyl 2,5,8,11,14-pentaoxa... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.057	2.48 ng/ul	192908	Naphthalene-d8	10.757

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Propanol, 1-[2-(2-methoxy-1-me...	206	C10H22O4	020324-33-8	86		
2	Tri(1,2-propyleneglycol), monome...	206	C10H22O4	1000262-26-6	74		
3	2-Propanol, 1-[2-(2-methoxy-1-me...	206	C10H22O4	020324-33-8	72		
4	Methyl 2,5,8,11,14-pentaoxahexad...	280	C12H24O7	1000366-81-5	72		
5	Methyl 2,5,8,11,14,17,20-heptaox...	368	C16H32O9	1010366-81-3	72		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102121\
Data File : BP007552.D
Acq On : 21 Oct 2021 13:29
Operator : CG/JU
Sample : M4262-04
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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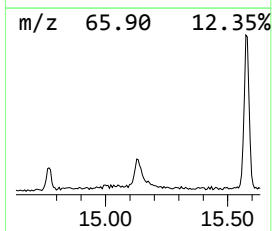
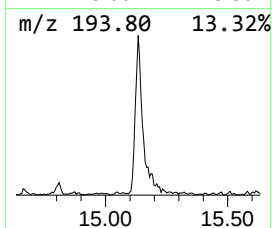
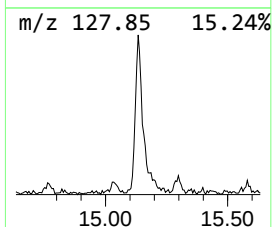
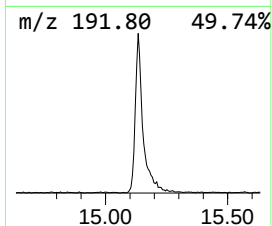
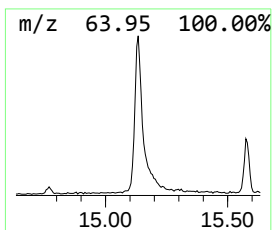
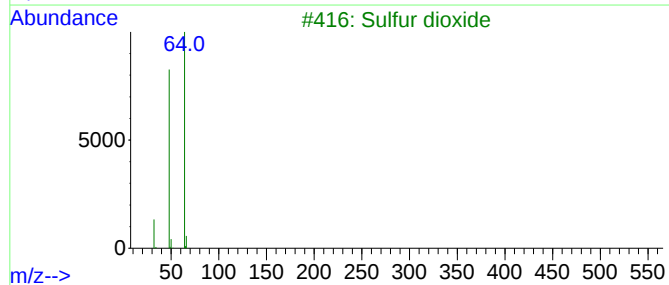
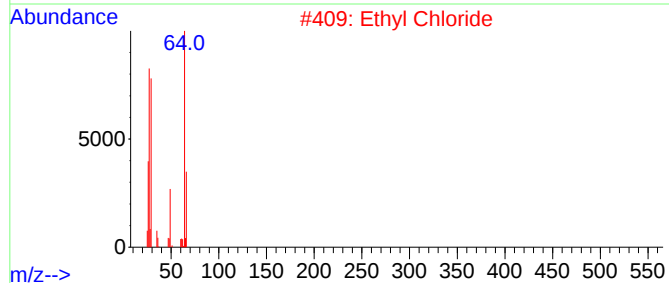
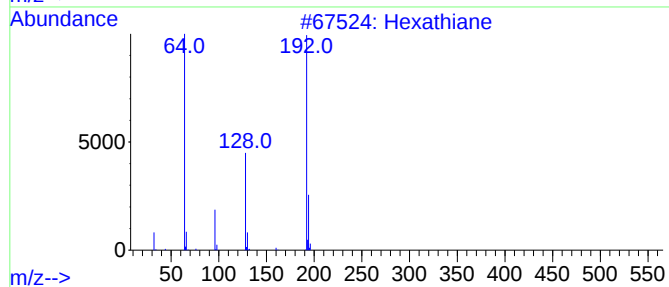
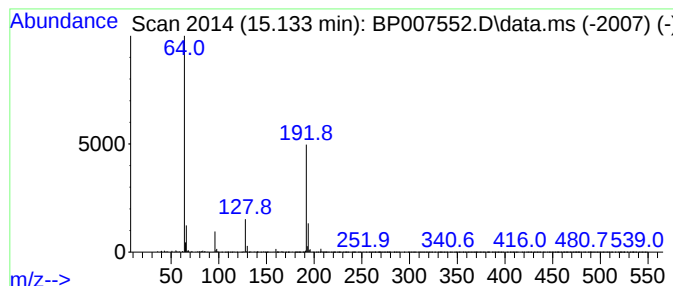
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Hexathiane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.133	2.25 ng/ul	237696	Acenaphthene-d10	14.586

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexathiane	192	S6	013798-23-7	91
2			Ethyl Chloride	64	C2H5Cl	000075-00-3	5
3			Sulfur dioxide	64	O2S	007446-09-5	4
4			Ethyl Chloride	64	C2H5Cl	000075-00-3	4
5			Ethyl Chloride	64	C2H5Cl	000075-00-3	4



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102121\
Data File : BP007552.D
Acq On : 21 Oct 2021 13:29
Operator : CG/JU
Sample : M4262-04
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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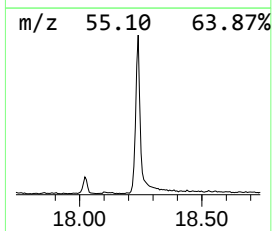
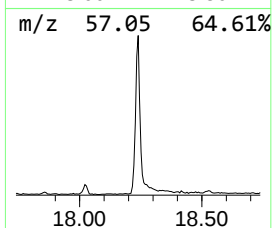
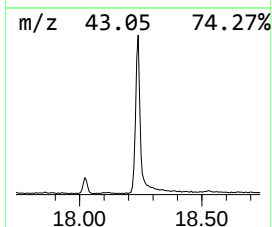
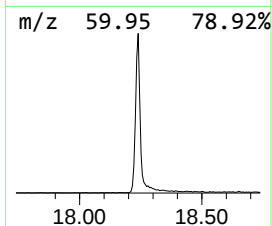
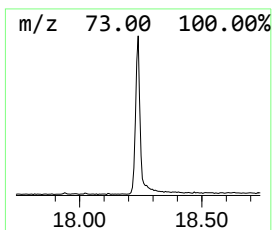
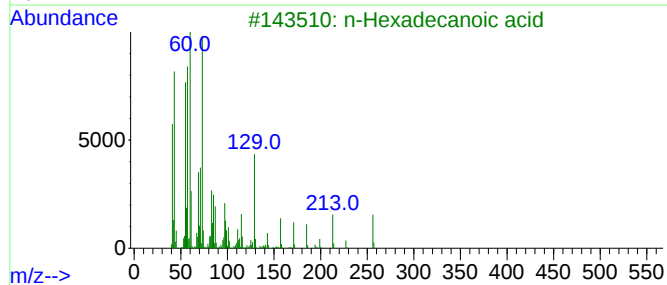
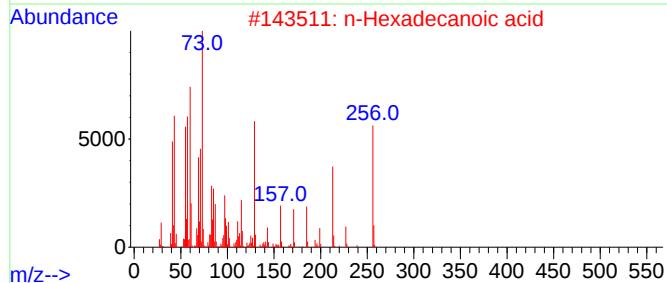
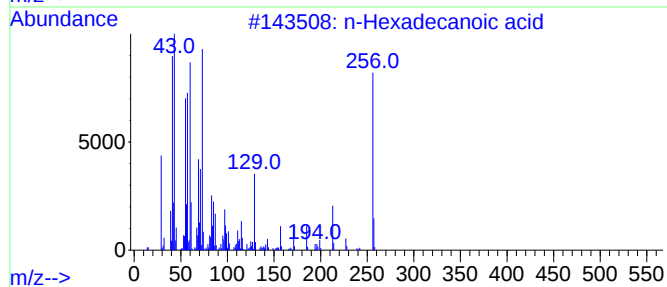
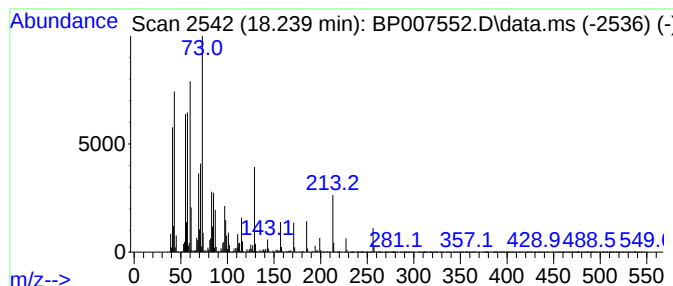
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 n-Hexadecanoic acid Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.239	12.17 ng/ul	1577940	Phenanthrene-d10	17.339

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
3			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
5			Tridecanoic acid	214	C13H26O2	000638-53-9	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102121\
Data File : BP007552.D
Acq On : 21 Oct 2021 13:29
Operator : CG/JU
Sample : M4262-04
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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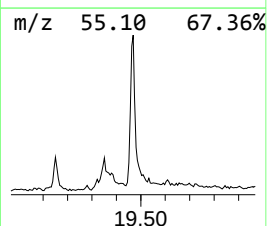
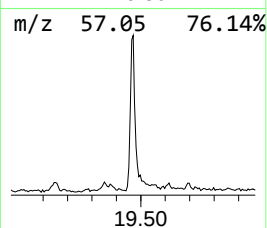
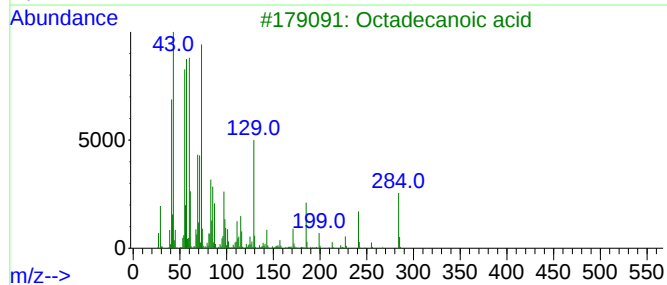
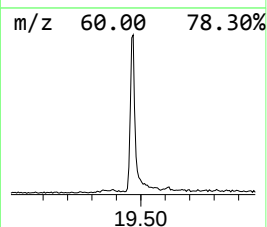
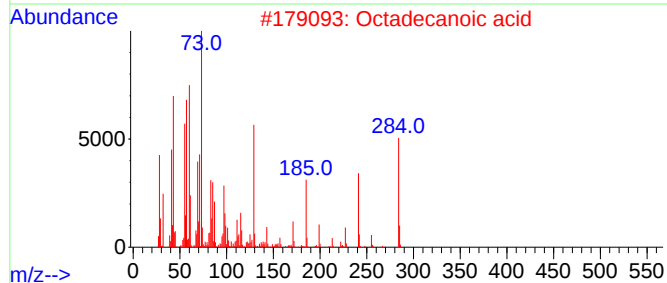
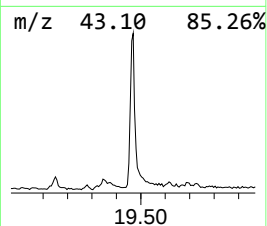
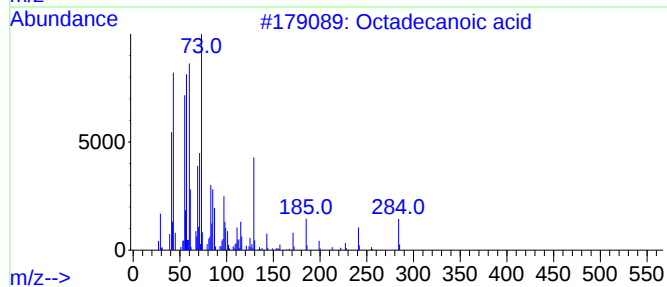
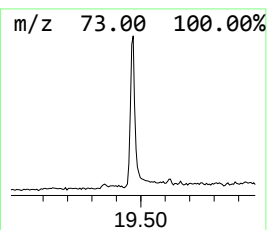
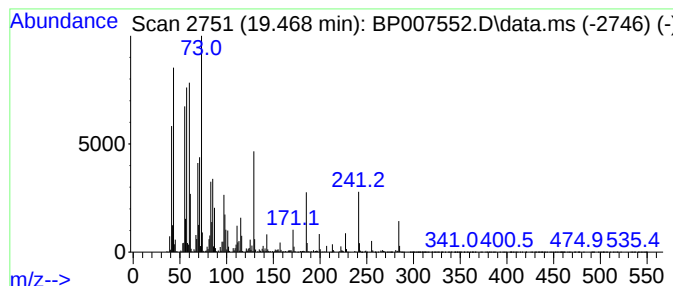
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Octadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.468	5.90 ng/ul	907079	Chrysene-d12	21.421

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102121\
Data File : BP007552.D
Acq On : 21 Oct 2021 13:29
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Sample : M4262-04
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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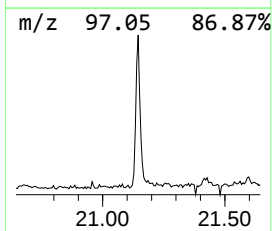
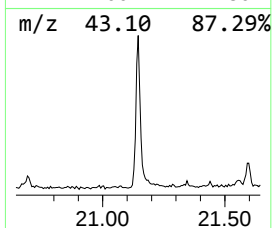
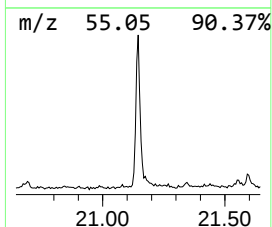
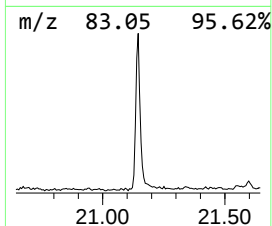
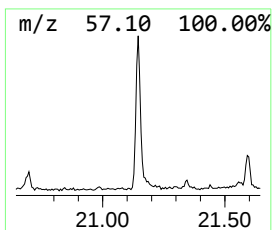
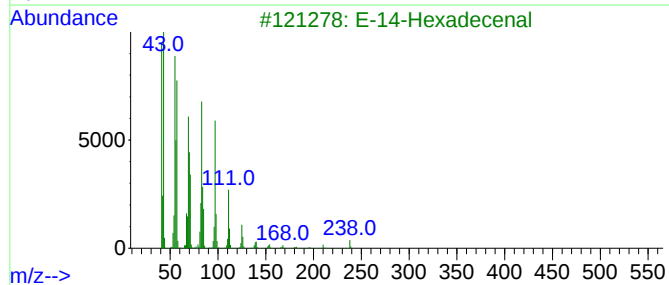
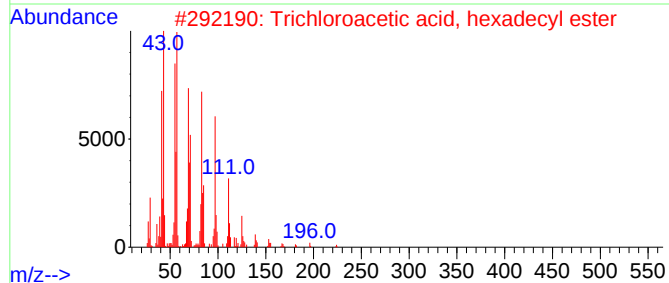
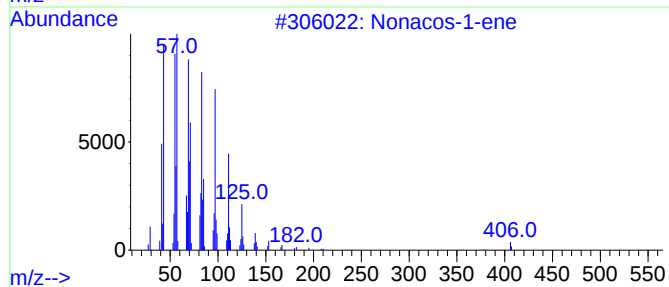
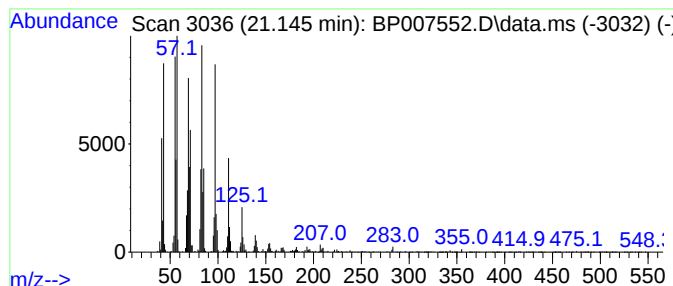
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.145	3.69 ng/ul	568133	Chrysene-d12	21.421

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonacos-1-ene	406	C29H58	018835-35-3	94
2			Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	94
3			E-14-Hexadecenal	238	C16H30O	330207-53-9	93
4			1-Heneicosanol	312	C21H44O	015594-90-8	93
5			1-Hexacosene	364	C26H52	018835-33-1	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102121\
 Data File : BP007552.D
 Acq On : 21 Oct 2021 13:29
 Operator : CG/JU
 Sample : M4262-04
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

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

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
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2-Propanol, 1-...	11.951	2.2	ng/ul	170389	2	10.757	1555650	20.0
Tri(1,2-propyle...	11.998	2.3	ng/ul	175000	2	10.757	1555650	20.0
Silane, trimeth...	12.016	2.5	ng/ul	190871	2	10.757	1555650	20.0
Methyl 2,5,8,11...	12.057	2.5	ng/ul	192908	2	10.757	1555650	20.0
Hexathiane	15.133	2.3	ng/ul	237696	3	14.586	2117490	20.0
n-Hexadecanoic ...	18.239	12.2	ng/ul	1577940	4	17.339	2593940	20.0
Octadecanoic acid	19.468	5.9	ng/ul	907079	5	21.421	3075580	20.0
(DEL) Alkane: S...	21.145	3.7	ng/ul	568133	5	21.421	3075580	20.0