

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP120423\
 Data File : BP018618.D
 Acq On : 05 Dec 2023 23:38
 Operator : MA/JU
 Sample : 05424-03
 Misc :
 ALS Vial : 54 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 YCGB7

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-BP120123.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP018618.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.264	27	30	37	rVB	66493	95748	0.79%	0.086%
2	3.558	75	80	83	rBV3	15140	26175	0.22%	0.023%
3	3.687	98	102	115	rBV	494083	741504	6.10%	0.662%
4	3.922	139	142	147	rVB2	19665	30128	0.25%	0.027%
5	4.016	154	158	162	rBV2	23922	31021	0.26%	0.028%
6	5.175	351	355	358	rBV5	9752	15765	0.13%	0.014%
7	7.016	658	668	678	rVB	409050	680537	5.60%	0.608%
8	7.187	690	697	706	rVB	1432300	2187601	18.00%	1.954%
9	7.381	723	730	739	rBV	1257117	1973258	16.24%	1.762%
10	7.463	740	744	751	rVB3	24893	47362	0.39%	0.042%
11	7.569	756	762	773	rVB	407976	660451	5.43%	0.590%
12	7.852	802	810	817	rBV	1231305	1911723	15.73%	1.707%
13	7.922	817	822	828	rVB	94480	158219	1.30%	0.141%
14	8.093	844	851	859	rBV	276521	455231	3.75%	0.407%
15	8.299	879	886	891	rBV2	74156	116231	0.96%	0.104%
16	8.357	891	896	905	rVB2	69385	112441	0.93%	0.100%
17	8.563	923	931	937	rBV	1367776	2146737	17.66%	1.917%
18	8.640	937	944	954	rVB2	95850	213779	1.76%	0.191%
19	9.010	998	1007	1014	rBV	1693462	2691773	22.15%	2.404%
20	9.075	1014	1018	1025	rVB	64836	108905	0.90%	0.097%
21	9.175	1032	1035	1040	rVB5	13438	20966	0.17%	0.019%
22	9.257	1043	1049	1054	rVB4	30288	51498	0.42%	0.046%
23	9.399	1066	1073	1077	rBV	55072	99360	0.82%	0.089%
24	9.734	1122	1130	1142	rBV	1237588	2043000	16.81%	1.825%
25	9.881	1148	1155	1161	rBV3	38237	80415	0.66%	0.072%
26	9.951	1161	1167	1175	rVB3	31113	76729	0.63%	0.069%
27	10.081	1181	1189	1197	rBV	359934	568095	4.67%	0.507%
28	10.269	1212	1221	1237	rBV	2101064	3544824	29.17%	3.166%
29	10.434	1242	1249	1257	rVV4	26929	55657	0.46%	0.050%
30	10.646	1277	1285	1297	rBV	1789162	3014721	24.80%	2.692%
31	10.787	1302	1309	1321	rBV	1686598	2791537	22.97%	2.493%
32	11.322	1394	1400	1403	rBV4	7853	12605	0.10%	0.011%
33	11.440	1415	1420	1423	rBV4	13008	17606	0.14%	0.016%
34	11.663	1454	1458	1462	rVB4	9077	13617	0.11%	0.012%
35	11.816	1476	1484	1491	rBV	53887	101210	0.83%	0.090%
36	11.887	1491	1496	1501	rBV5	29232	64639	0.53%	0.058%

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120123.MA.M
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37	12.145	1535	1540	1546	rBV3	15814	27149	0.22%	0.024%
38	12.234	1549	1555	1558	rBV	44519	75696	0.62%	0.068%
39	12.269	1558	1561	1566	rVB2	21572	33014	0.27%	0.029%
40	12.440	1586	1590	1595	rVV2	17896	25729	0.21%	0.023%

41	12.851	1654	1660	1665	rBV	107994	162339	1.34%	0.145%
42	12.940	1672	1675	1680	rBV5	11662	16783	0.14%	0.015%
43	13.075	1692	1698	1700	rBV2	13751	24692	0.20%	0.022%
44	13.275	1728	1732	1733	rBV2	10879	15387	0.13%	0.014%
45	13.387	1745	1751	1759	rVB2	72778	117053	0.96%	0.105%

46	13.451	1759	1762	1764	rBV3	10580	14172	0.12%	0.013%
47	13.704	1804	1805	1810	rVB4	10884	12470	0.10%	0.011%
48	13.898	1831	1838	1846	rBV	4583658	6381978	52.51%	5.699%
49	14.010	1855	1857	1863	rVB5	9505	14524	0.12%	0.013%
50	14.175	1875	1885	1896	rBV	5032242	7221770	59.42%	6.449%

51	14.269	1897	1901	1909	rVB4	24568	51834	0.43%	0.046%
52	14.351	1911	1915	1919	rBV5	26037	36957	0.30%	0.033%
53	14.481	1926	1937	1948	rBV	3076918	4562745	37.54%	4.075%
54	14.563	1948	1951	1955	rVB6	17276	25440	0.21%	0.023%
55	14.681	1965	1971	1980	rBV	327257	512423	4.22%	0.458%

56	14.834	1992	1997	2002	rVB6	18857	34748	0.29%	0.031%
57	14.898	2002	2008	2011	rBV7	27932	48432	0.40%	0.043%
58	14.945	2011	2016	2021	rVV2	22895	42215	0.35%	0.038%
59	15.051	2029	2034	2043	rVB	320377	453192	3.73%	0.405%
60	15.228	2060	2064	2069	rBV8	11122	13445	0.11%	0.012%

61	15.281	2069	2073	2075	rBV	16430	20294	0.17%	0.018%
62	15.328	2075	2081	2085	rVB8	22051	50131	0.41%	0.045%
63	15.475	2099	2106	2117	rBV	6479161	9645936	79.37%	8.614%
64	15.586	2118	2125	2137	rBV	1853893	2637332	21.70%	2.355%
65	15.710	2142	2146	2148	rBV2	29250	39609	0.33%	0.035%

66	15.745	2148	2152	2157	rVB2	109094	169556	1.40%	0.151%
67	15.804	2158	2162	2165	rBV2	29956	39218	0.32%	0.035%
68	15.875	2166	2174	2180	rBV2	386590	727296	5.98%	0.650%
69	15.969	2187	2190	2197	rVB4	49275	85211	0.70%	0.076%
70	16.116	2211	2215	2220	rVB5	34842	44976	0.37%	0.040%

71	16.210	2225	2231	2244	rVV	741446	1321913	10.88%	1.181%
72	16.322	2248	2250	2254	rVV3	24418	28952	0.24%	0.026%
73	16.369	2254	2258	2264	rVB	105881	140341	1.15%	0.125%
74	16.522	2282	2284	2295	rVB	19801	35787	0.29%	0.032%
75	16.628	2297	2302	2307	rBV5	39297	66253	0.55%	0.059%

76	16.716	2314	2317	2321	rBV	34977	45564	0.37%	0.041%
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 Title : SVOA CALIBRATION

77	16.980	2358	2362	2366	rBV5	34357	63649	0.52%	0.057%
78	17.022	2366	2369	2379	rVB3	66670	118913	0.98%	0.106%
79	17.110	2381	2384	2389	rVB6	14496	24384	0.20%	0.022%
80	17.228	2397	2404	2413	rBV	3745544	5531445	45.51%	4.940%
81	17.328	2414	2421	2429	rVV	7482859	10582166	87.07%	9.450%
82	17.904	2516	2519	2523	rVB3	19402	23731	0.20%	0.021%
83	18.122	2550	2556	2565	rBV	925341	1297629	10.68%	1.159%
84	18.433	2606	2609	2612	rBV3	16656	22475	0.18%	0.020%
85	18.539	2622	2627	2631	rBV	153006	205580	1.69%	0.184%
86	18.927	2690	2693	2697	rBV	50228	58813	0.48%	0.053%
87	19.092	2718	2721	2724	rVB2	33472	36940	0.30%	0.033%
88	19.139	2725	2729	2735	rVB	109907	141273	1.16%	0.126%
89	19.204	2736	2740	2744	rBV	115565	153603	1.26%	0.137%
90	19.351	2761	2765	2774	rBV	364642	516588	4.25%	0.461%
91	19.492	2786	2789	2794	rVB4	39555	50334	0.41%	0.045%
92	19.563	2795	2801	2808	rBV	9136135	12153842	100.00%	10.854%
93	21.033	3047	3051	3059	rBV2	377810	688968	5.67%	0.615%
94	21.310	3093	3098	3104	rBV	3752112	4900405	40.32%	4.376%
95	23.527	3468	3475	3483	rBV2	4965457	9098148	74.86%	8.125%
96	23.680	3494	3501	3511	rVB	2160553	4322388	35.56%	3.860%

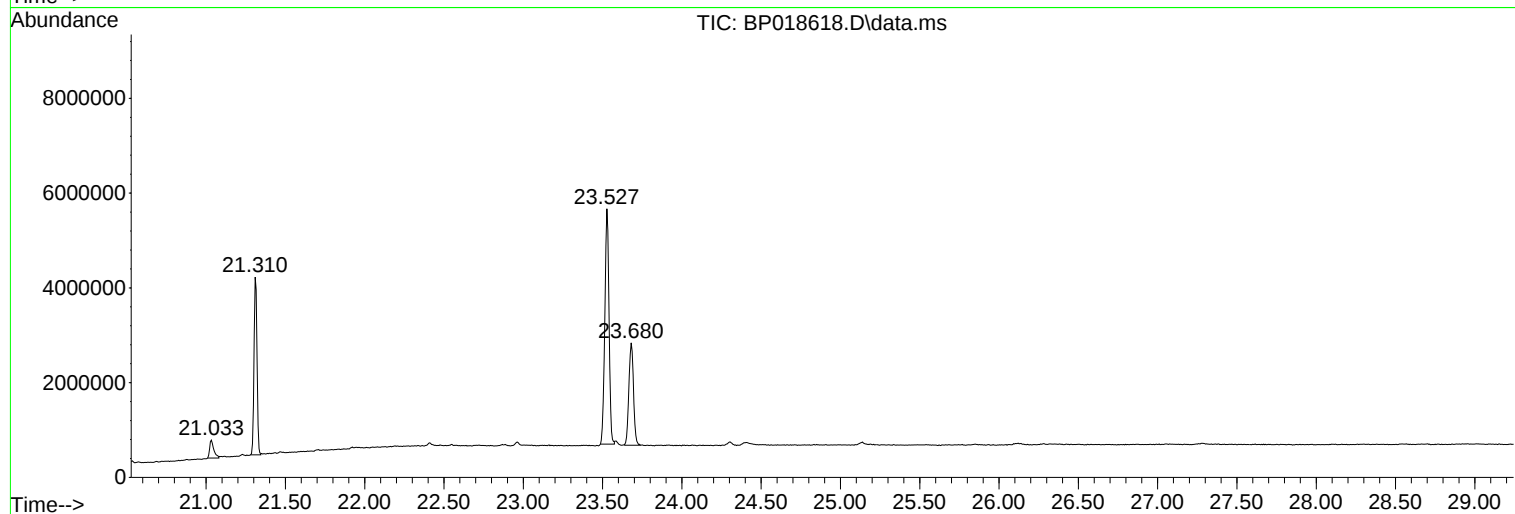
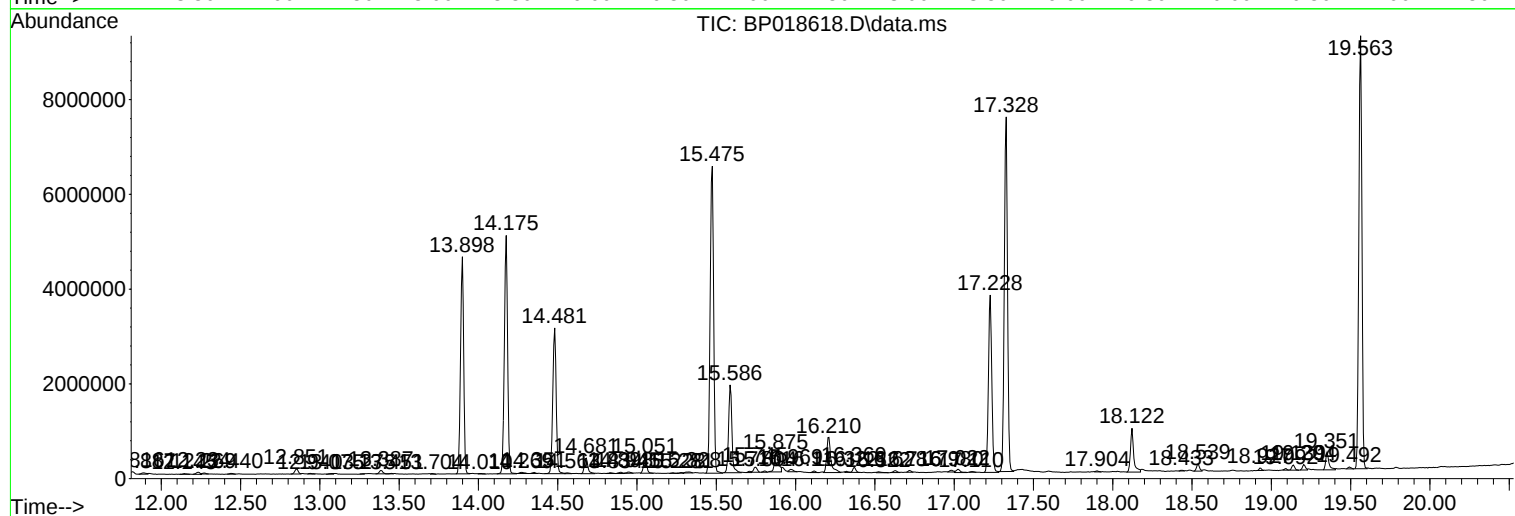
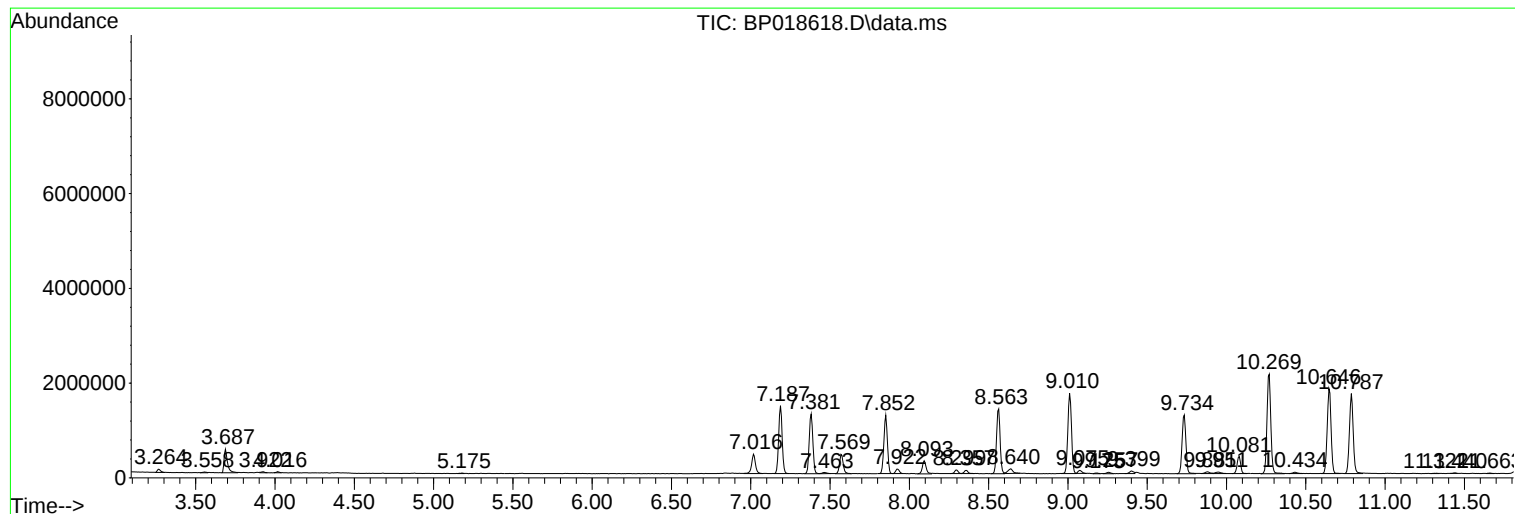
Sum of corrected areas: 111974898

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120123.MA.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP120423\
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ALS Vial : 54 Sample Multiplier: 1

Instrument :
BNA_P
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YCGB7

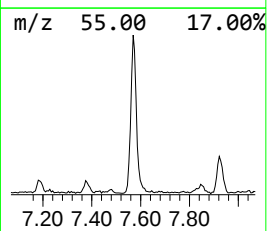
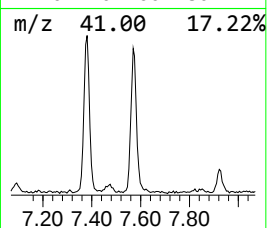
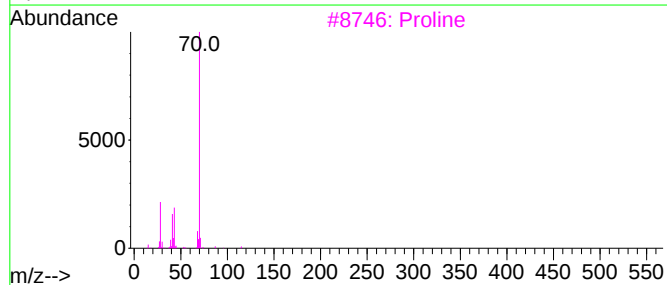
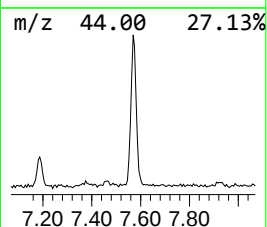
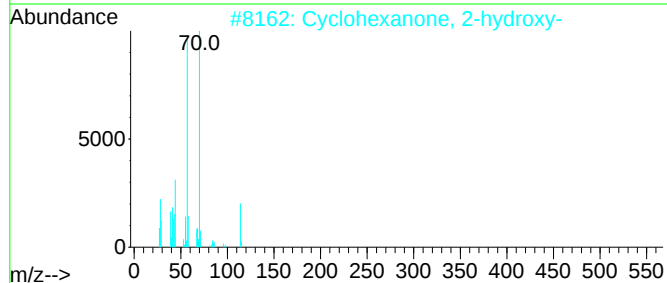
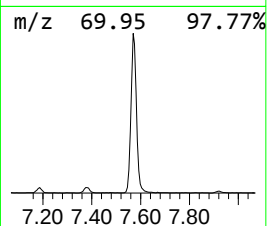
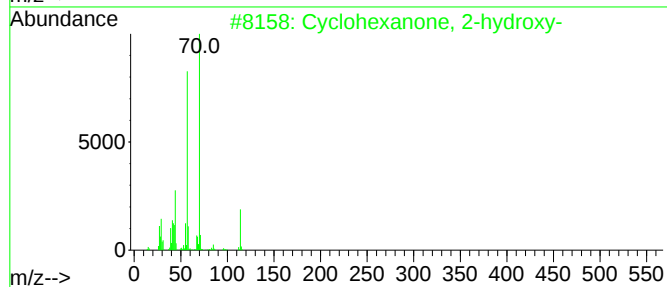
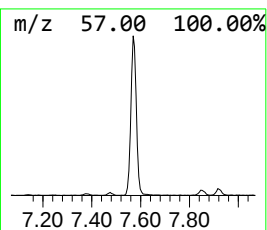
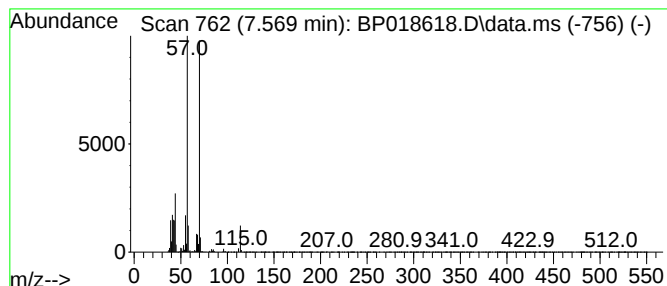
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120123.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Cyclohexanone, 2-hydroxy- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.569	6.91 ng/ul	660451	1,4-Dichlorobenzene-d4	7.852

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexanone, 2-hydroxy-	114	C6H10O2	000533-60-8	83
2			Cyclohexanone, 2-hydroxy-	114	C6H10O2	000533-60-8	58
3			Proline	115	C5H9NO2	000147-85-3	43
4			D-Proline	115	C5H9NO2	000344-25-2	43
5			dl-Proline	115	C5H9NO2	000609-36-9	37



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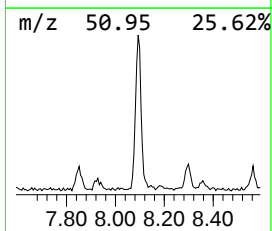
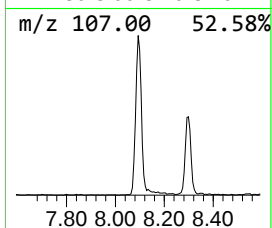
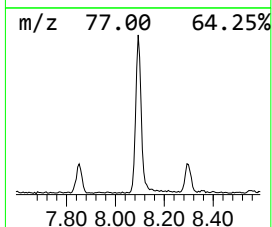
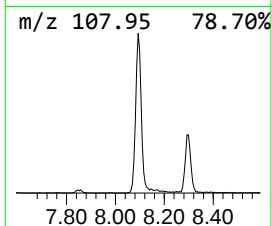
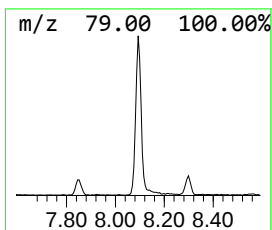
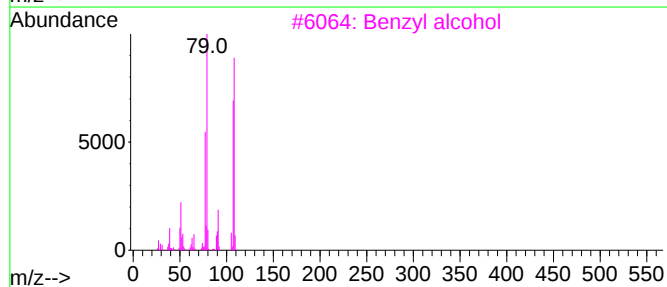
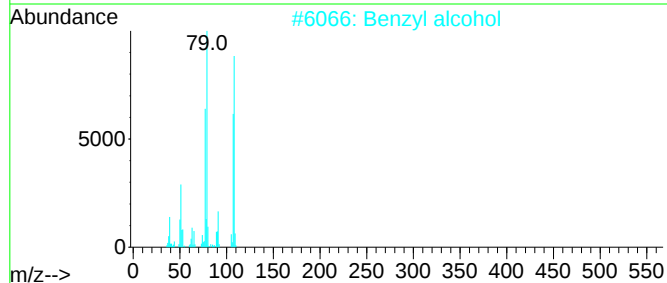
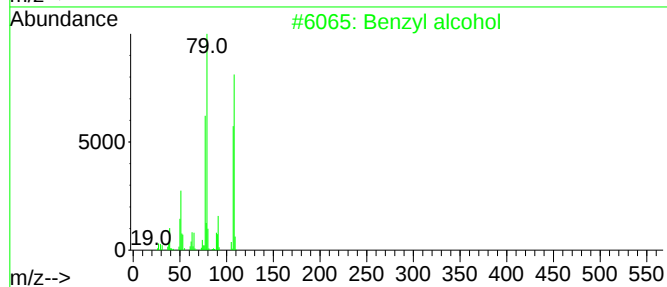
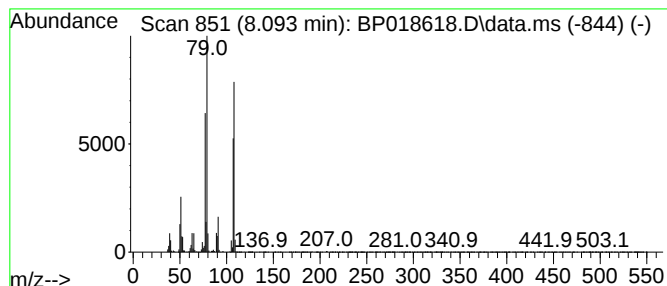
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120123.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Benzyl alcohol Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.093	4.76 ng/ul	455231	1,4-Dichlorobenzene-d4	7.852

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzyl alcohol	108	C7H8O	000100-51-6	98
2			Benzyl alcohol	108	C7H8O	000100-51-6	98
3			Benzyl alcohol	108	C7H8O	000100-51-6	97
4			Benzyl alcohol	108	C7H8O	000100-51-6	95
5			N-Cbz-glycylglycine p-nitropheny...	387	C18H17N3O7	013574-81-7	91



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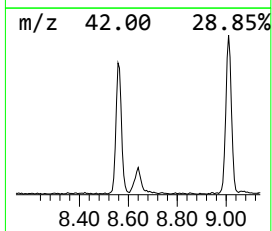
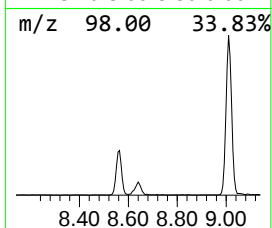
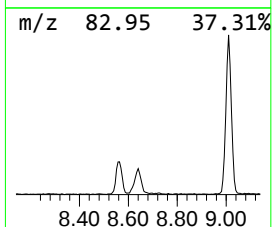
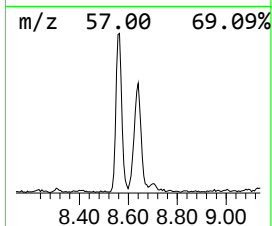
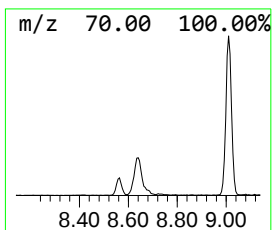
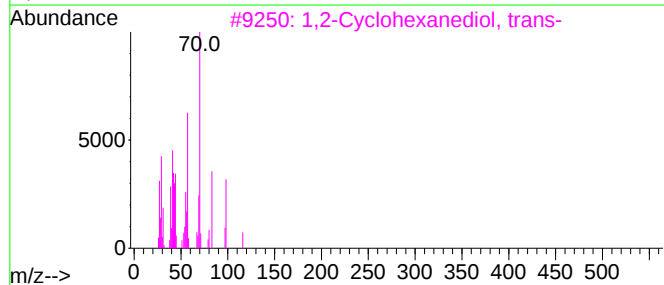
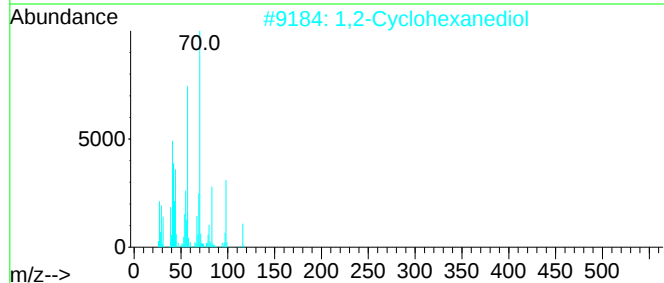
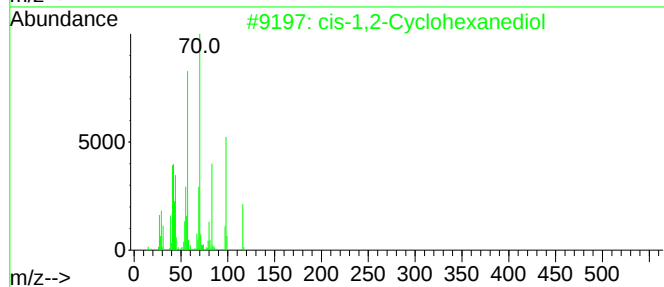
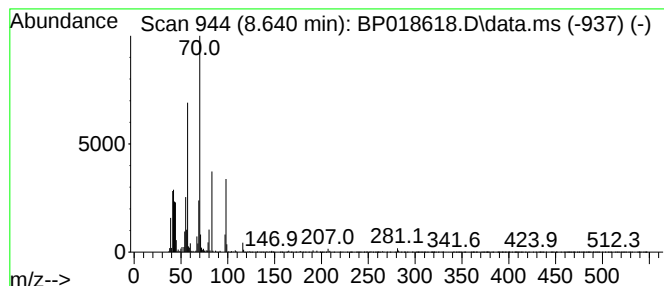
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 cis-1,2-Cyclohexanediol Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.640	2.24 ng/ul	213779	1,4-Dichlorobenzene-d4	7.852

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			cis-1,2-Cyclohexanediol	116	C6H12O2	001792-81-0	93
2			1,2-Cyclohexanediol	116	C6H12O2	000931-17-9	91
3			1,2-Cyclohexanediol, trans-	116	C6H12O2	001460-57-7	90
4			cis-1,2-Cyclohexanediol	116	C6H12O2	001792-81-0	90
5			1,2-Cyclohexanediol	116	C6H12O2	000931-17-9	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP120423\
Data File : BP018618.D
Acq On : 05 Dec 2023 23:38
Operator : MA/JU
Sample : 05424-03
Misc :
ALS Vial : 54 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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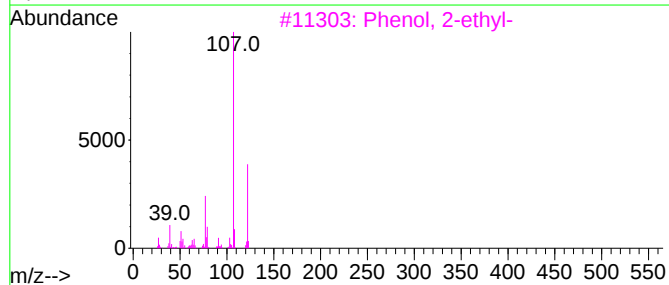
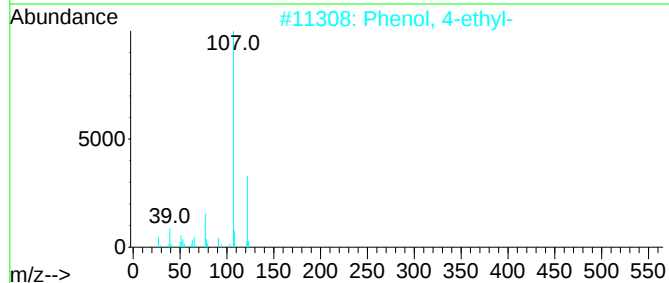
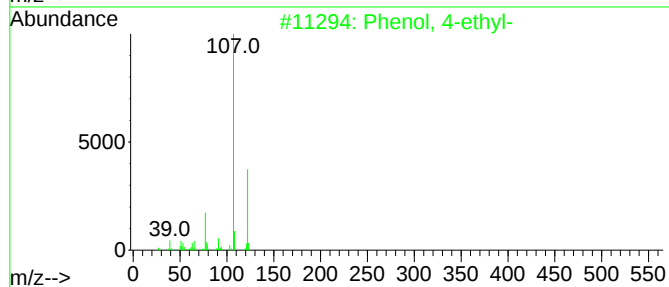
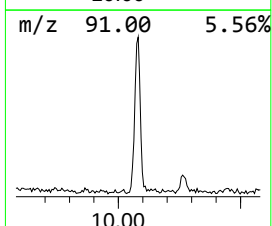
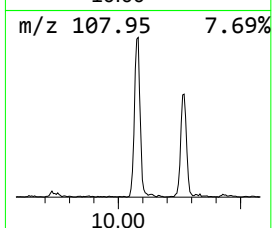
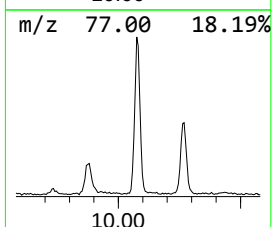
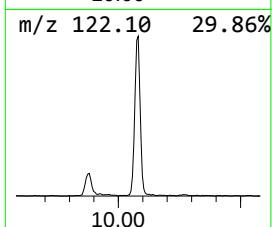
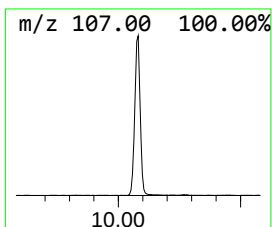
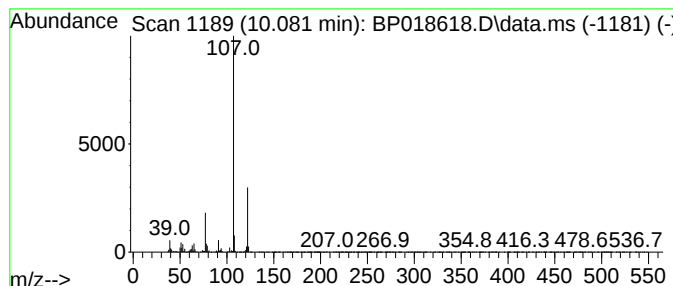
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Phenol, 4-ethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.081	3.77 ng/ul	568095	Naphthalene-d8	10.646

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenol, 4-ethyl-	122	C8H10O	000123-07-9	95
2			Phenol, 4-ethyl-	122	C8H10O	000123-07-9	95
3			Phenol, 2-ethyl-	122	C8H10O	000090-00-6	91
4			Phenol, 4-ethyl-	122	C8H10O	000123-07-9	91
5			Phenol, 2-ethyl-	122	C8H10O	000090-00-6	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP120423\
Data File : BP018618.D
Acq On : 05 Dec 2023 23:38
Operator : MA/JU
Sample : 05424-03
Misc :
ALS Vial : 54 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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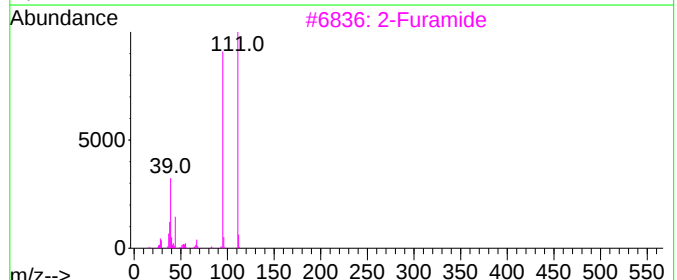
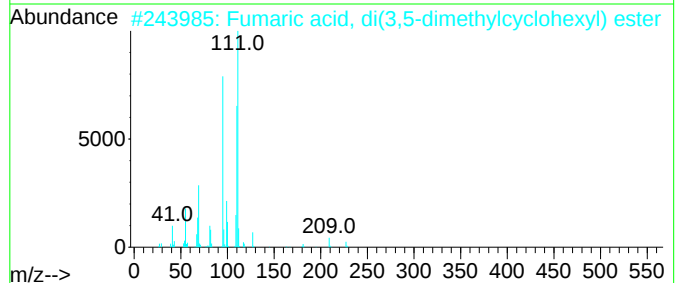
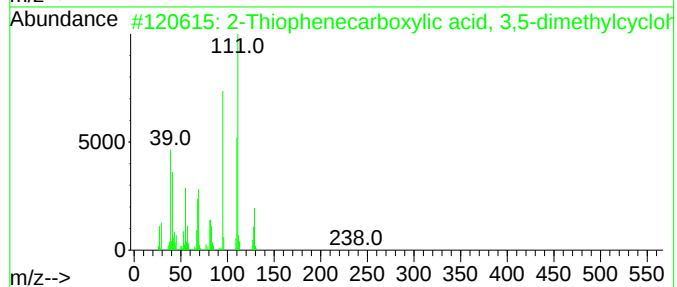
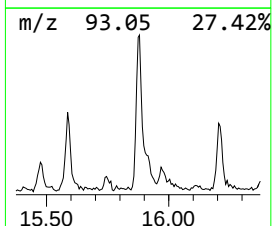
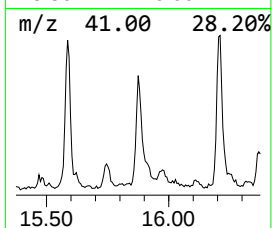
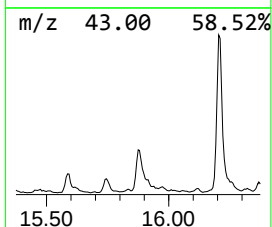
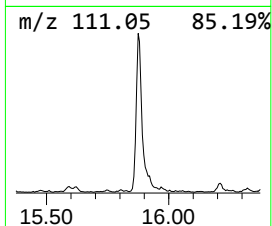
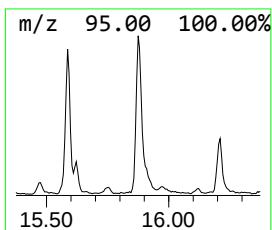
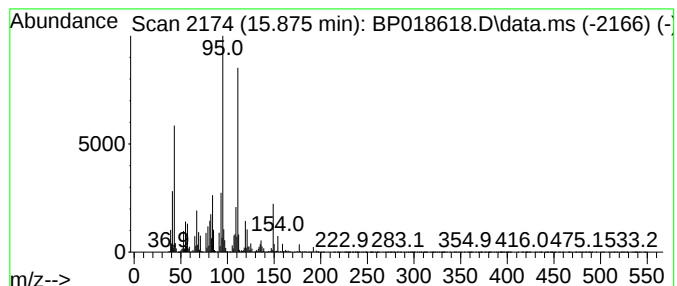
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 2-Thiophenecarboxylic acid,... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.875	2.63 ng/ul	727296	Phenanthrene-d10	17.227

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Thiophenecarboxylic acid, 3,5-...	238	C13H18O2S	1000278-87-5	50
2			Fumaric acid, di(3,5-dimethylcyc...	336	C20H32O4	1000345-06-5	50
3			2-Furamide	111	C5H5N02	000609-38-1	47
4			N-Hydroxy-4-pyridone	111	C5H5N02	1010426-43-6	46
5			4-Pyridinol-1-oxide	111	C5H5N02	006890-62-6	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP120423\
Data File : BP018618.D
Acq On : 05 Dec 2023 23:38
Operator : MA/JU
Sample : 05424-03
Misc :
ALS Vial : 54 Sample Multiplier: 1

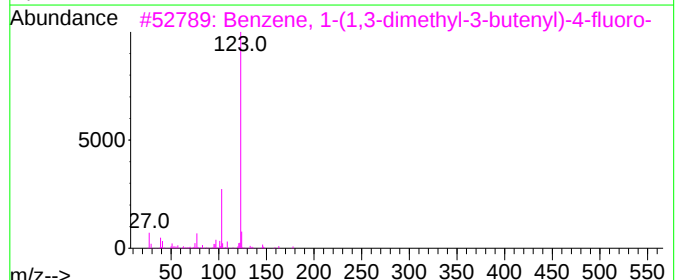
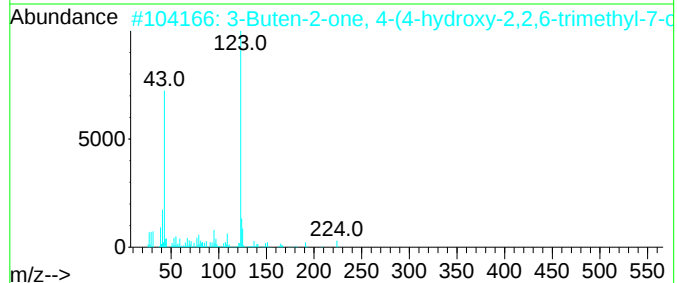
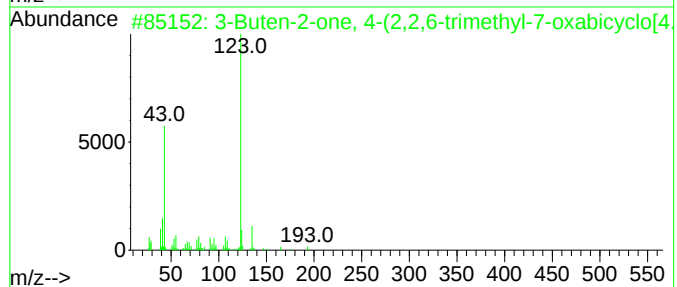
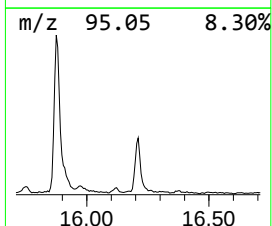
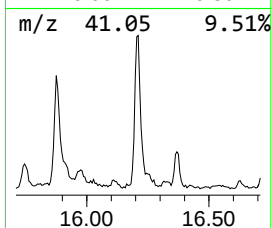
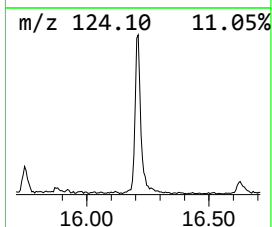
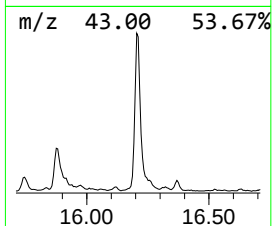
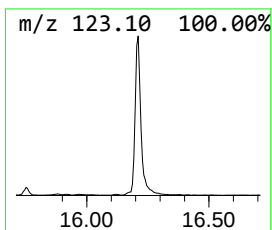
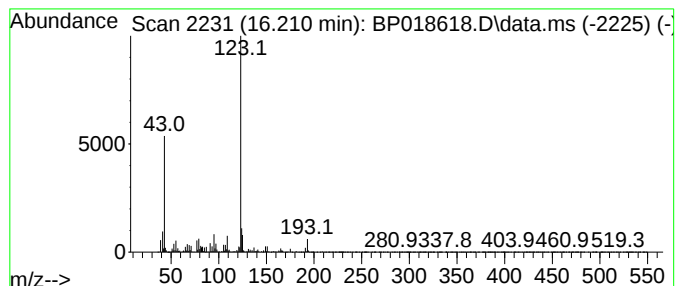
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120123.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 6 3-Buten-2-one, 4-(2,2,6-tri... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
16.210	4.78 ng/ul	1321910	Phenanthrene-d10	17.227	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Buten-2-one, 4-(2,2,6-trimethyl-3-buten-2-yl)benzene	208	C13H20O2	023267-57-4	87
2	3-Buten-2-one, 4-(4-hydroxy-2,2,6-trimethyl-3-buten-2-yl)benzene	224	C13H20O3	038274-01-0	83
3	Benzene, 1-(1,3-dimethyl-3-buten-2-yl)-	178	C12H15F	074646-34-7	64
4	Cyclopropanecarboxylic acid, 2,2,4,4-tetramethyl-	300	C19H24O3	023031-36-9	64
5	3-Fluorobenzoic acid, tridec-2-ynyl ester	318	C20H27FO2	1000292-60-9	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP120423\
Data File : BP018618.D
Acq On : 05 Dec 2023 23:38
Operator : MA/JU
Sample : 05424-03
Misc :
ALS Vial : 54 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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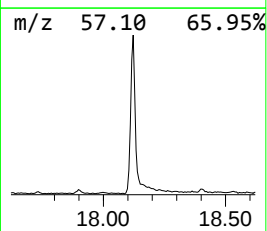
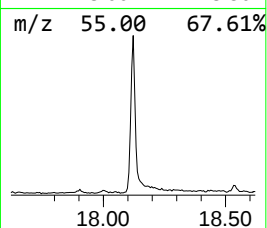
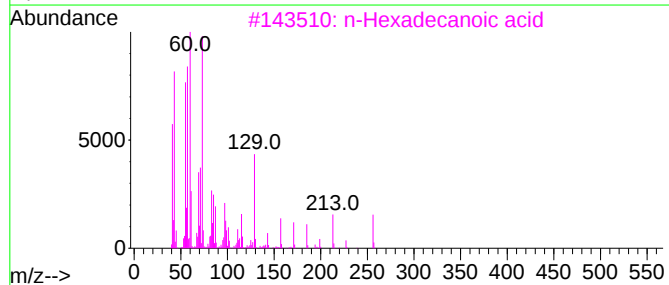
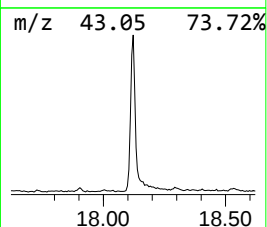
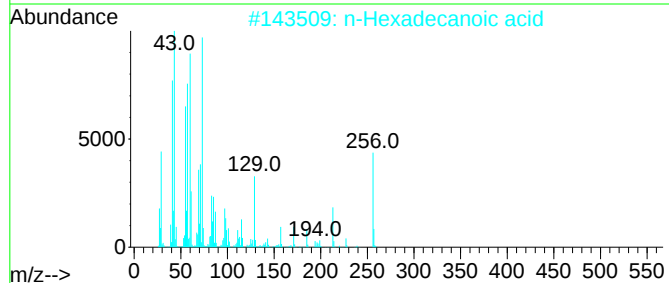
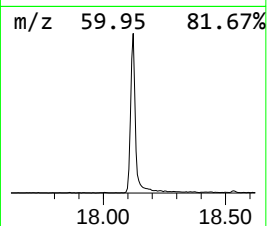
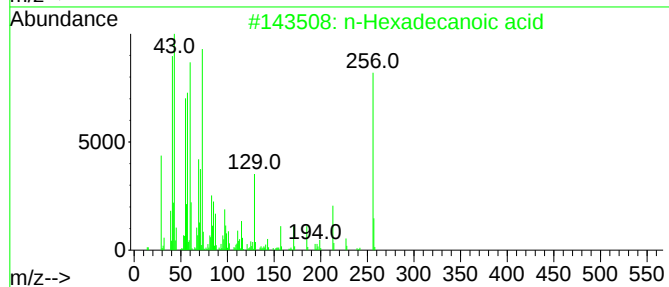
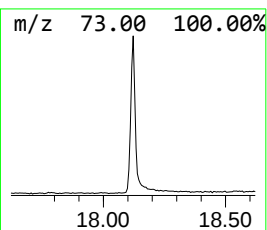
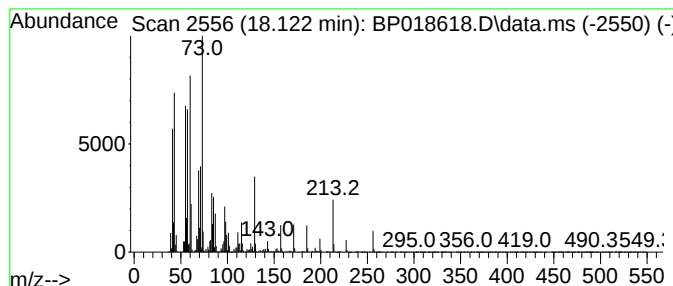
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 n-Hexadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.122	4.69 ng/ul	1297630	Phenanthrene-d10	17.227

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			Tridecanoic acid	214	C13H26O2	000638-53-9	91
5			Tetradecanoic acid	228	C14H28O2	000544-63-8	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP120423\
Data File : BP018618.D
Acq On : 05 Dec 2023 23:38
Operator : MA/JU
Sample : 05424-03
Misc :
ALS Vial : 54 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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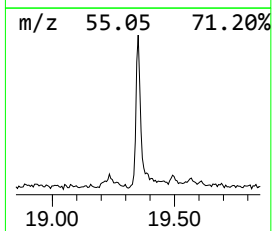
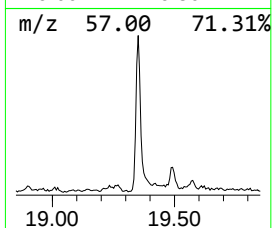
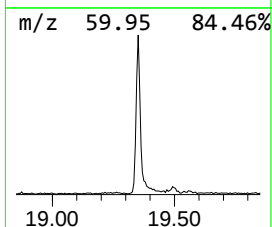
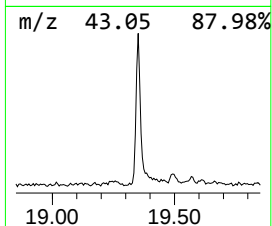
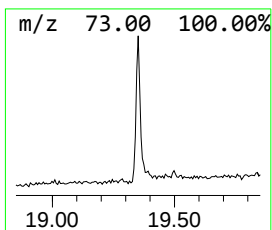
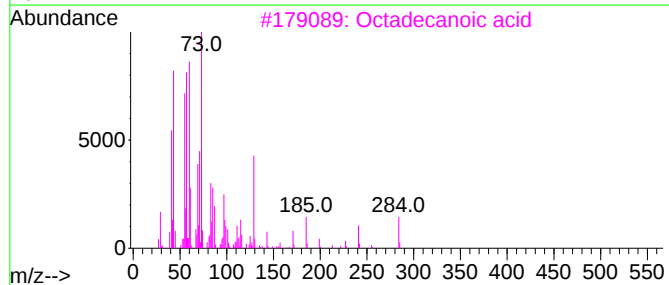
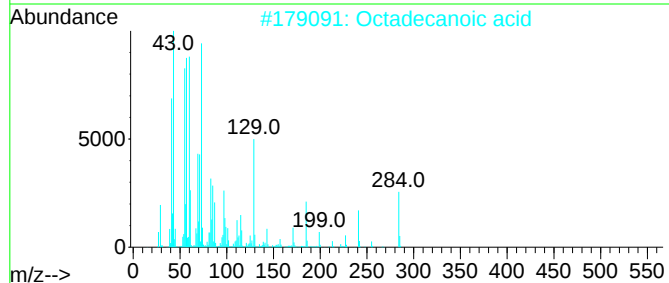
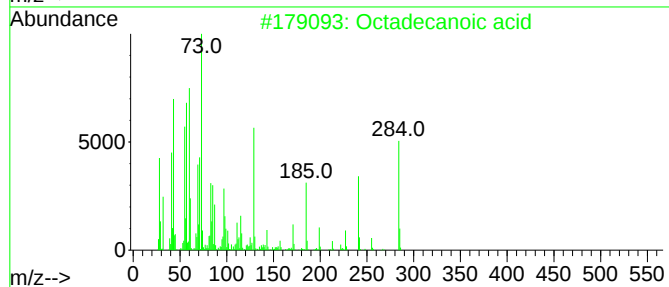
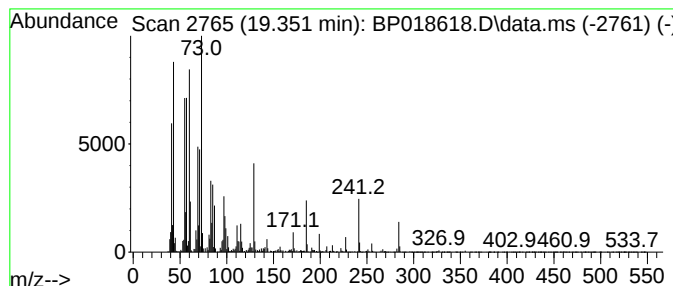
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Octadecanoic acid Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.351	2.11 ng/ul	516588	Chrysene-d12	21.310

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	94
5			Pentadecanoic acid	242	C15H30O2	001002-84-2	89



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP120423\
Data File : BP018618.D
Acq On : 05 Dec 2023 23:38
Operator : MA/JU
Sample : 05424-03
Misc :
ALS Vial : 54 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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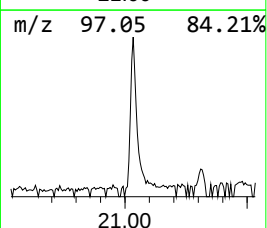
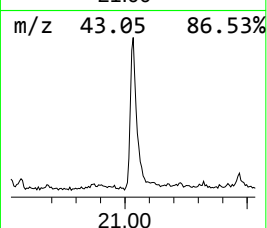
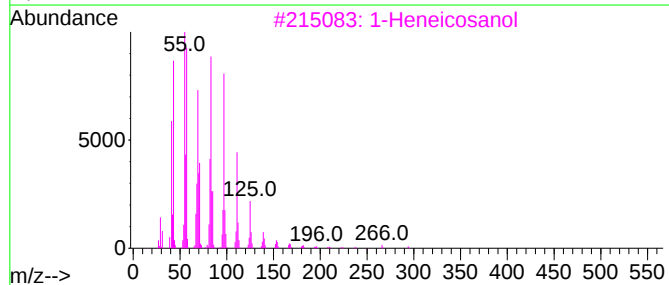
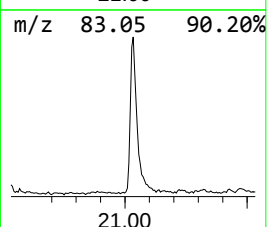
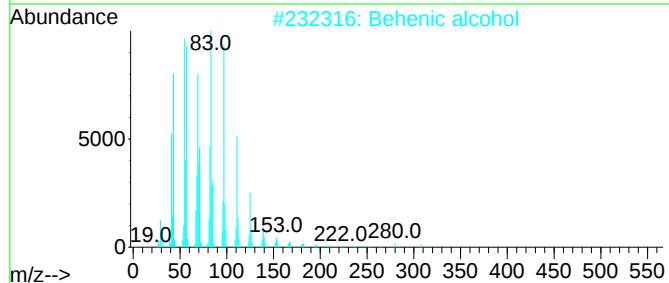
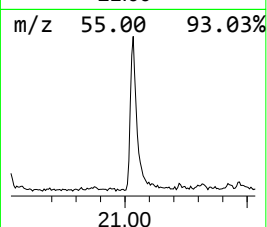
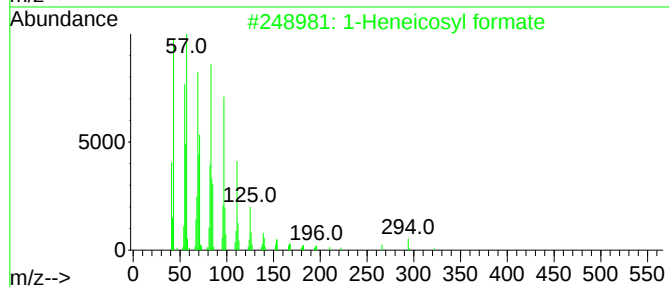
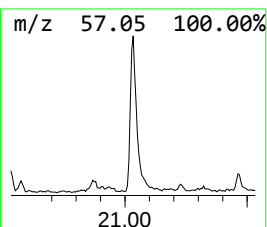
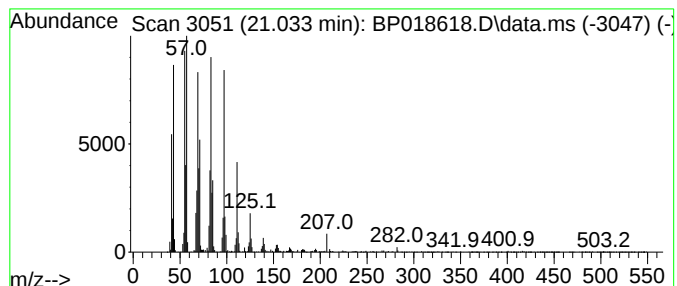
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 1-Heneicosyl formate Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.033	2.81 ng/ul	688968	Chrysene-d12	21.310

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Heneicosyl formate	340	C22H44O2	077899-03-7	95
2			Behenic alcohol	326	C22H46O	000661-19-8	94
3			1-Heneicosanol	312	C21H44O	015594-90-8	94
4			Pentacos-1-ene	350	C25H50	016980-85-1	94
5			n-Tetracosanol-1	354	C24H50O	000506-51-4	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP120423\
Data File : BP018618.D
Acq On : 05 Dec 2023 23:38
Operator : MA/JU
Sample : 05424-03
Misc :
ALS Vial : 54 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120123.MA.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
Cyclohexanone, ...	7.569	6.9	ng/ul	660451	1	7.852	1911720	20.0
Benzyl alcohol	8.093	4.8	ng/ul	455231	1	7.852	1911720	20.0
cis-1,2-Cyclohe...	8.640	2.2	ng/ul	213779	1	7.852	1911720	20.0
Phenol, 4-ethyl-	10.081	3.8	ng/ul	568095	2	10.646	3014720	20.0
2-Thiophenecarb...	15.875	2.6	ng/ul	727296	4	17.227	5531450	20.0
3-Buten-2-one, ...	16.210	4.8	ng/ul	1321910	4	17.227	5531450	20.0
n-Hexadecanoic ...	18.122	4.7	ng/ul	1297630	4	17.227	5531450	20.0
Octadecanoic acid	19.351	2.1	ng/ul	516588	5	21.310	4900410	20.0
1-Heneicosyl fo...	21.033	2.8	ng/ul	688968	5	21.310	4900410	20.0