

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072324\
 Data File : BP021114.D
 Acq On : 24 Jul 2024 01:51
 Operator : MA/JU
 Sample : P3238-06
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 A4BE9

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-BP071024.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP021114.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.193	31	35	42	rVB	27826	43240	0.75%	0.079%
2	3.481	80	84	100	rVB	63371	111920	1.94%	0.203%
3	3.605	102	105	106	rVV3	7679	7944	0.14%	0.014%
4	3.622	106	108	117	rVV	59658	103573	1.79%	0.188%
5	3.705	118	122	128	rVB	41449	59930	1.04%	0.109%
6	3.911	154	157	162	rVB	22288	30638	0.53%	0.056%
7	4.069	180	184	190	rVB9	3552	7691	0.13%	0.014%
8	4.240	210	213	217	rBV6	4101	5905	0.10%	0.011%
9	4.322	224	227	231	rBV	5301	8271	0.14%	0.015%
10	4.422	238	244	248	rBV7	5993	9113	0.16%	0.017%
11	4.463	248	251	255	rVV4	11083	14047	0.24%	0.026%
12	4.816	304	311	320	rVB	40083	68966	1.19%	0.125%
13	5.781	472	475	481	rBV2	5459	9431	0.16%	0.017%
14	6.763	636	642	647	rBV4	11073	22339	0.39%	0.041%
15	6.875	655	661	678	rBV	407058	773746	13.40%	1.407%
16	7.010	678	684	696	rVB	355583	572408	9.92%	1.041%
17	7.210	711	718	727	rBV	497869	804502	13.94%	1.463%
18	7.299	727	733	740	rVV2	21787	62799	1.09%	0.114%
19	7.663	786	795	804	rBV	544838	913273	15.82%	1.661%
20	7.746	804	809	815	rVB9	6272	13606	0.24%	0.025%
21	8.399	913	920	932	rBV	503866	911356	15.79%	1.657%
22	8.499	933	937	944	rVB3	11504	24240	0.42%	0.044%
23	8.822	985	992	1007	rBV	440034	772466	13.38%	1.405%
24	8.969	1012	1017	1022	rVB8	7979	16531	0.29%	0.030%
25	9.169	1045	1051	1055	rBV5	9461	15337	0.27%	0.028%
26	9.369	1080	1085	1098	rVB2	36576	74167	1.28%	0.135%
27	9.534	1106	1113	1126	rBV	393846	674229	11.68%	1.226%
28	9.810	1157	1160	1164	rBV6	6141	10741	0.19%	0.020%
29	10.081	1198	1206	1225	rBV	520457	1062174	18.40%	1.931%
30	10.428	1258	1265	1271	rBV	788909	1318289	22.84%	2.397%
31	10.475	1271	1273	1280	rVB	29552	42130	0.73%	0.077%
32	10.593	1285	1293	1312	rBV	321239	701199	12.15%	1.275%
33	10.981	1351	1359	1368	rBV8	17487	43360	0.75%	0.079%
34	11.187	1387	1394	1403	rBV	118378	263174	4.56%	0.479%
35	11.940	1518	1522	1526	rBV6	3379	6256	0.11%	0.011%
36	12.028	1531	1537	1543	rVB3	9863	17504	0.30%	0.032%

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
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37	12.098	1543	1549	1555	rBV	16798	28869	0.50%	0.052%
38	12.322	1582	1587	1591	rBV2	11671	19085	0.33%	0.035%
39	12.751	1652	1660	1661	rBV8	3397	6192	0.11%	0.011%
40	13.087	1713	1717	1719	rBV5	5289	6109	0.11%	0.011%
41	13.251	1740	1745	1752	rBV10	5563	11804	0.20%	0.021%
42	13.481	1773	1784	1785	rBV7	7881	23186	0.40%	0.042%
43	13.616	1799	1807	1813	rBV4	13276	28404	0.49%	0.052%
44	13.716	1813	1824	1830	rBV	984329	1510320	26.16%	2.746%
45	13.857	1844	1848	1851	rVB5	5649	7027	0.12%	0.013%
46	13.992	1860	1871	1882	rBV	1241385	1992027	34.51%	3.622%
47	14.175	1899	1902	1908	rVB8	6060	10647	0.18%	0.019%
48	14.298	1915	1923	1930	rBV	1181909	1855399	32.14%	3.374%
49	14.369	1930	1935	1947	rVB	109726	176500	3.06%	0.321%
50	14.516	1952	1960	1969	rVB5	72829	194030	3.36%	0.353%
51	14.604	1969	1975	1989	rBV	259659	652058	11.30%	1.186%
52	14.716	1990	1994	1998	rBV2	64556	95501	1.65%	0.174%
53	14.816	2009	2011	2019	rVB9	10009	14623	0.25%	0.027%
54	14.934	2026	2031	2036	rBV5	8076	16161	0.28%	0.029%
55	15.098	2053	2059	2063	rBV7	15144	35975	0.62%	0.065%
56	15.322	2090	2097	2103	rBV	1502876	2424401	42.00%	4.408%
57	15.381	2103	2107	2113	rVB2	122499	171148	2.96%	0.311%
58	15.481	2118	2124	2130	rBV	250056	378078	6.55%	0.687%
59	15.675	2153	2157	2164	rVB	22797	33605	0.58%	0.061%
60	15.798	2175	2178	2179	rBV3	7224	7967	0.14%	0.014%
61	15.828	2179	2183	2189	rVV2	28888	51057	0.88%	0.093%
62	15.881	2189	2192	2195	rBV4	18879	23588	0.41%	0.043%
63	16.039	2215	2219	2222	rBV3	16949	29514	0.51%	0.054%
64	16.210	2244	2248	2252	rBV6	12119	23173	0.40%	0.042%
65	16.286	2259	2261	2264	rBV3	8465	9656	0.17%	0.018%
66	16.628	2315	2319	2325	rBV6	33008	67578	1.17%	0.123%
67	16.769	2339	2343	2351	rVB	52344	89315	1.55%	0.162%
68	16.863	2352	2359	2364	rBV5	37156	93581	1.62%	0.170%
69	16.933	2366	2371	2376	rBV	71988	113366	1.96%	0.206%
70	16.986	2376	2380	2383	rBV	45342	60723	1.05%	0.110%
71	17.028	2383	2387	2393	rVB3	59690	92518	1.60%	0.168%
72	17.116	2396	2402	2406	rBV	1605631	2387650	41.36%	4.341%
73	17.163	2406	2410	2415	rVV	1543539	2223592	38.52%	4.043%
74	17.222	2415	2420	2425	rVV2	1648477	2646641	45.85%	4.812%
75	17.257	2425	2426	2431	rVB	233087	230814	4.00%	0.420%
76	17.304	2431	2434	2437	rBV2	36426	42599	0.74%	0.077%

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77	17.551	2467	2476	2480	rBV2	119002	289600	5.02%	0.527%
78	17.739	2501	2508	2515	rBV3	145626	324619	5.62%	0.590%
79	17.804	2516	2519	2522	rVB	46797	54836	0.95%	0.100%
80	18.051	2558	2561	2564	rBV	312246	331759	5.75%	0.603%
81	18.163	2577	2580	2585	rVB	68028	95839	1.66%	0.174%
82	18.227	2585	2591	2595	rBV3	312457	586719	10.16%	1.067%
83	18.263	2595	2597	2605	rVB2	125329	223955	3.88%	0.407%
84	18.327	2605	2608	2616	rBV3	66298	139288	2.41%	0.253%
85	18.404	2617	2621	2625	rBV2	84372	146138	2.53%	0.266%
86	18.539	2641	2644	2651	rVV	95644	164576	2.85%	0.299%
87	18.604	2651	2655	2662	rVB	209155	345818	5.99%	0.629%
88	19.257	2761	2766	2771	rVB	2958215	3708107	64.24%	6.742%
89	19.380	2784	2787	2797	rVB3	214688	367502	6.37%	0.668%
90	19.610	2820	2826	2829	rBV2	2679312	3984574	69.03%	7.245%
91	19.639	2829	2831	2838	rVB	2042126	2240604	38.82%	4.074%
92	20.151	2916	2918	2924	rVB	358556	430575	7.46%	0.783%
93	20.263	2934	2937	2941	rVB	168438	201046	3.48%	0.366%
94	21.139	3082	3086	3089	rBV2	220365	352755	6.11%	0.641%
95	21.216	3096	3099	3108	rVB4	585243	912338	15.80%	1.659%
96	21.563	3152	3158	3163	rBV3	3129536	5772490	100.00%	10.496%
97	21.610	3163	3166	3173	rVB	954789	1393956	24.15%	2.535%
98	23.845	3540	3546	3552	rBV2	441542	1155580	20.02%	2.101%
99	24.668	3679	3686	3692	rBV2	797818	1996696	34.59%	3.630%
100	24.892	3718	3724	3732	rVB2	864250	2297833	39.81%	4.178%

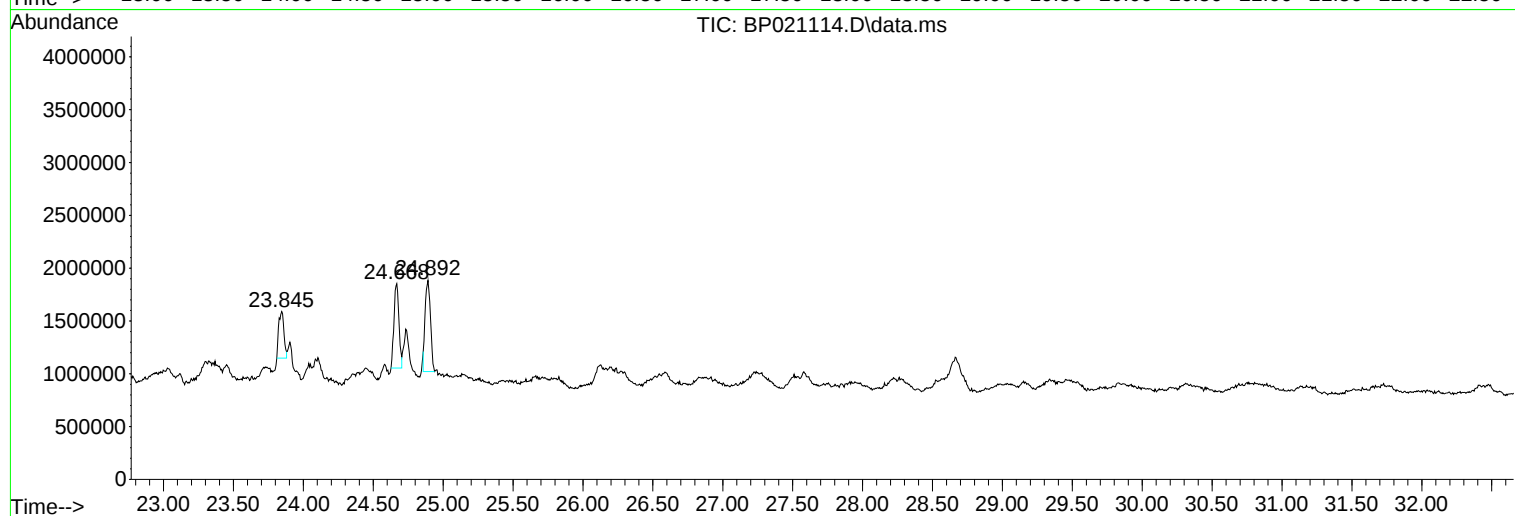
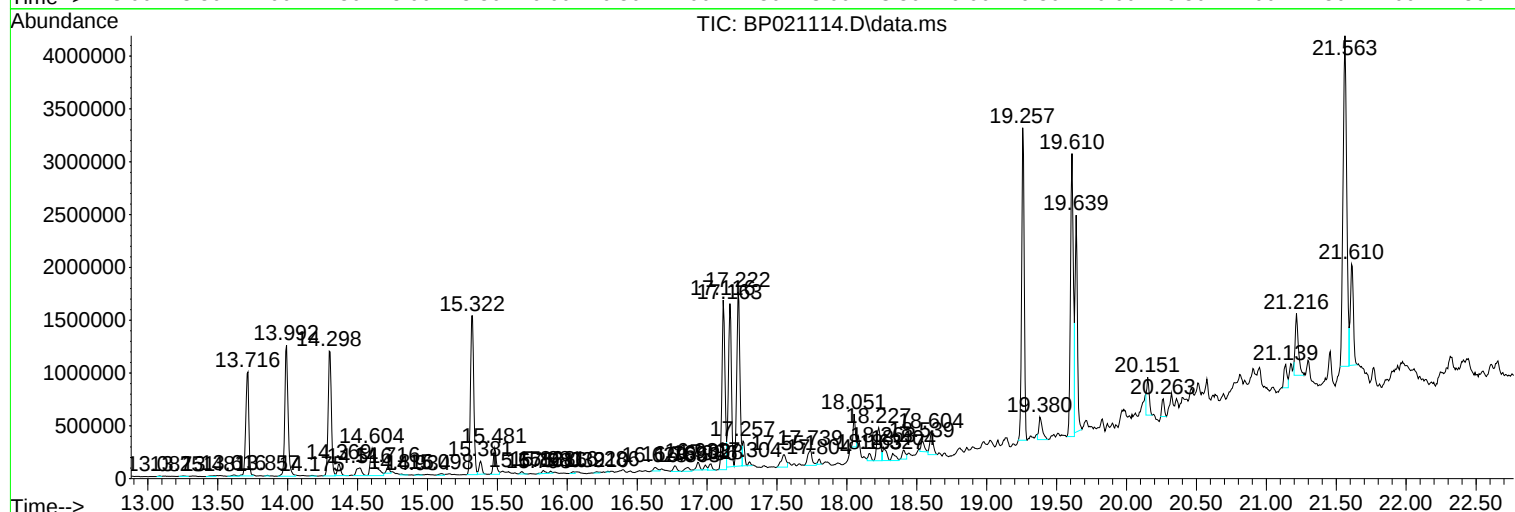
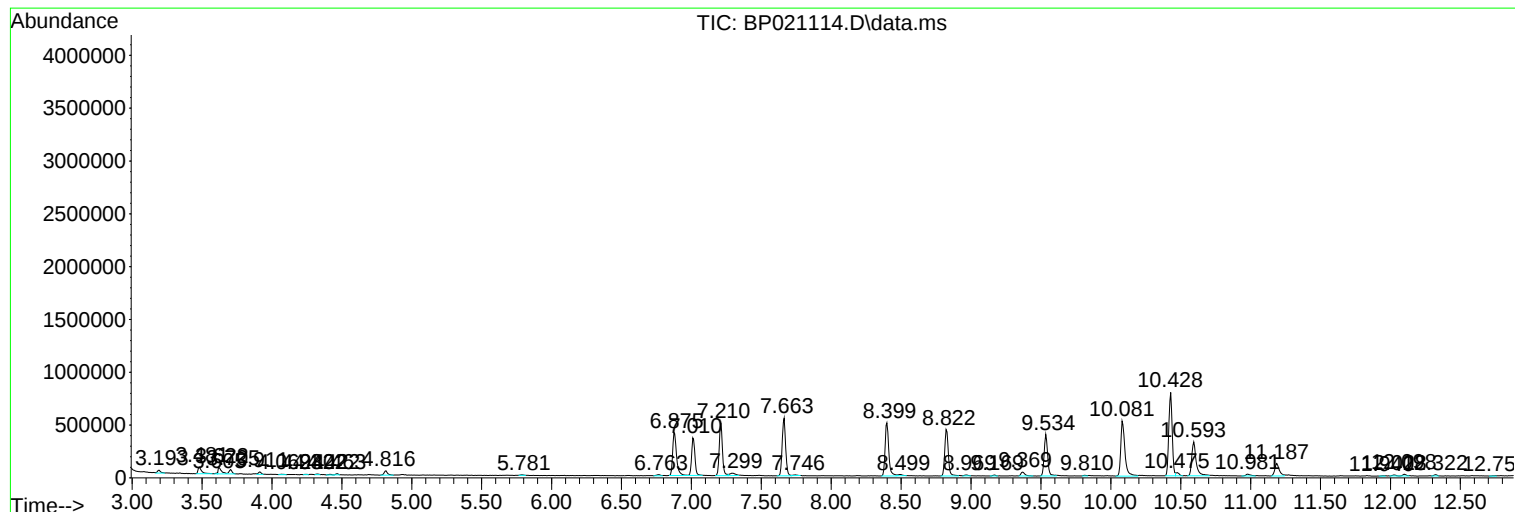
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

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



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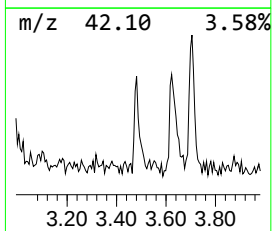
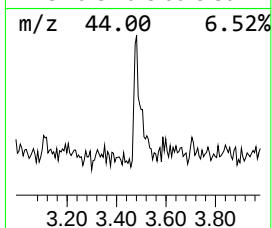
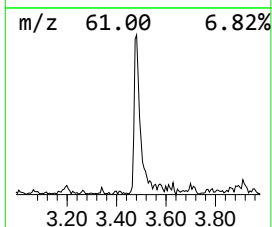
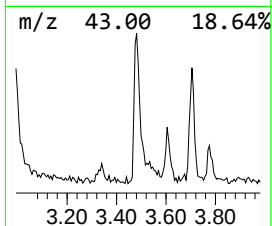
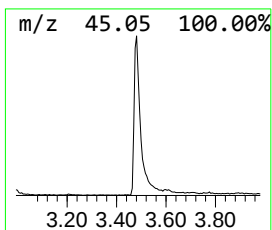
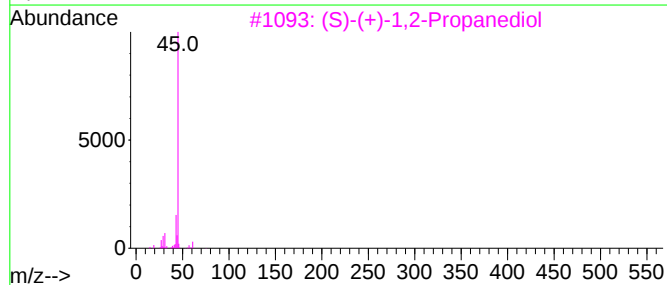
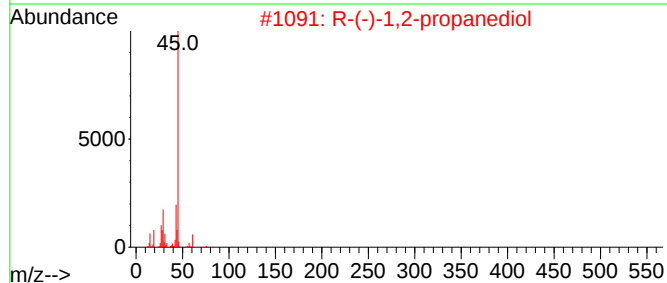
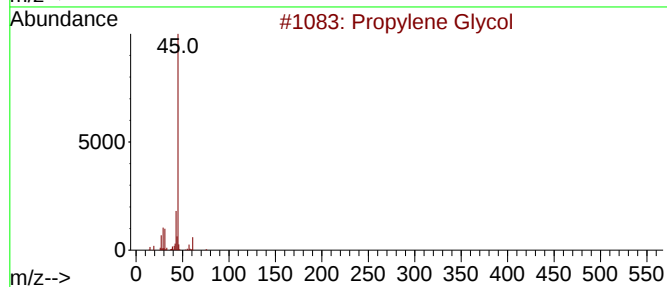
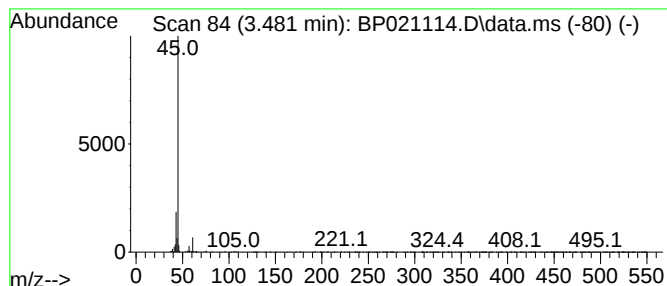
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Propylene Glycol Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.481	2.45 ng/ul	111920	1,4-Dichlorobenzene-d4	7.663

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Propylene Glycol	76	C3H8O2	000057-55-6	78
2			R-(-)-1,2-propanediol	76	C3H8O2	004254-14-2	72
3			(S)-(+)-1,2-Propanediol	76	C3H8O2	004254-15-3	40
4			Propanoic acid, 2-hydroxy-, meth...	104	C4H8O3	000547-64-8	40
5			Paraldehyde	132	C6H12O3	000123-63-7	9



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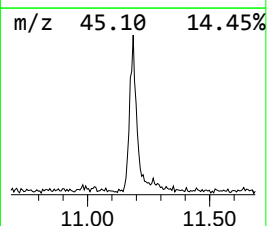
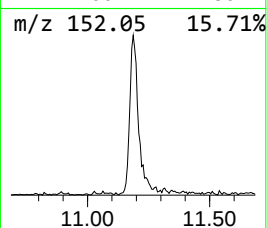
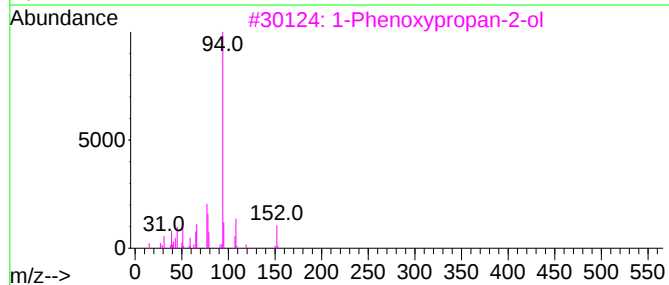
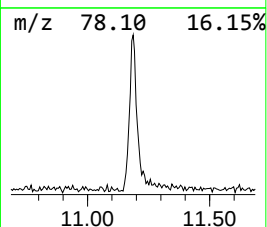
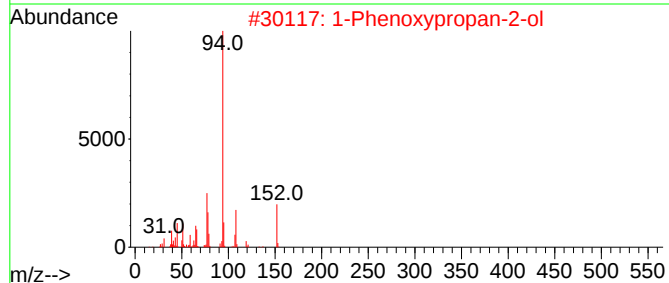
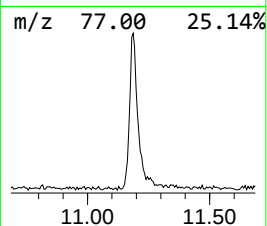
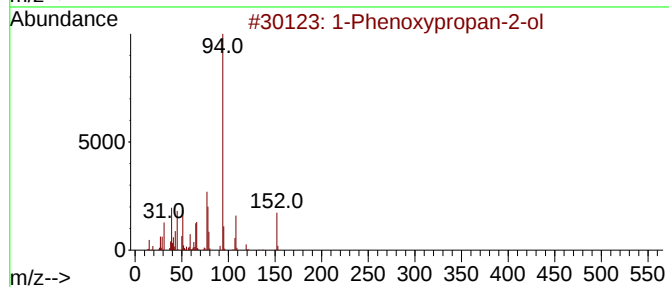
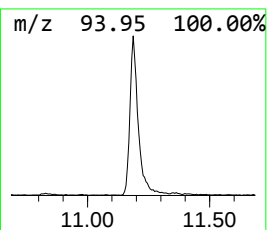
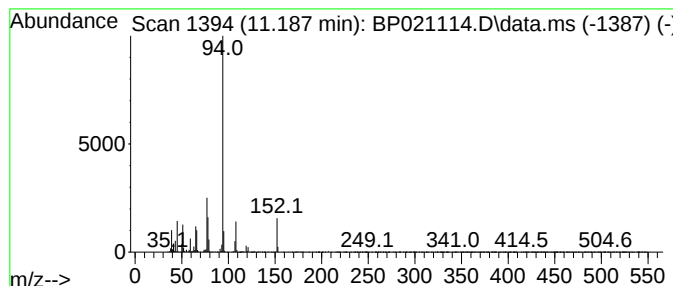
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 1-Phenoxypropan-2-ol Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.187	3.99 ng/ul	263174	Naphthalene-d8	10.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	96
2			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	96
3			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	95
4			1-Propanol, 3-phenoxy-	152	C9H12O2	006180-61-6	93
5			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	87



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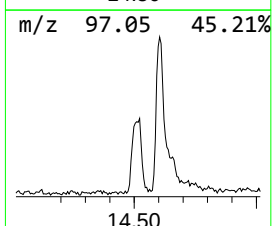
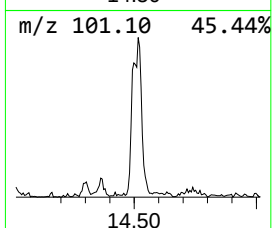
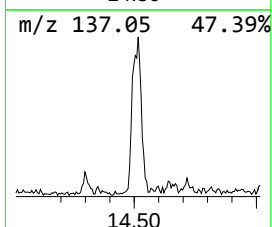
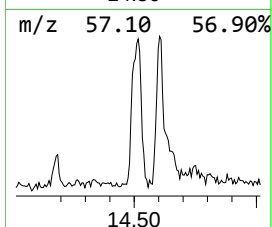
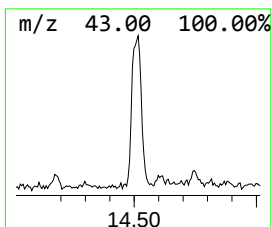
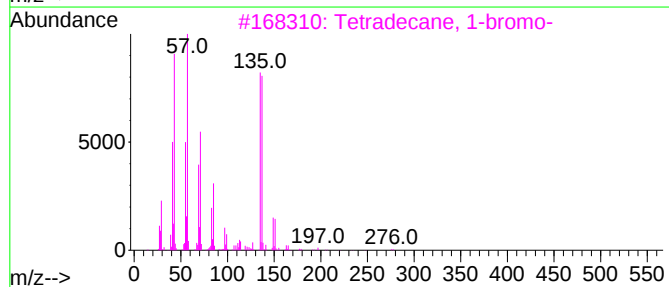
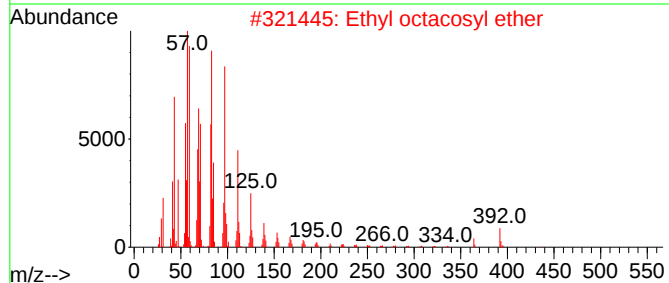
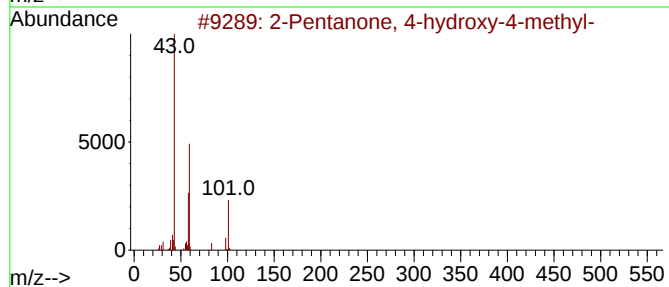
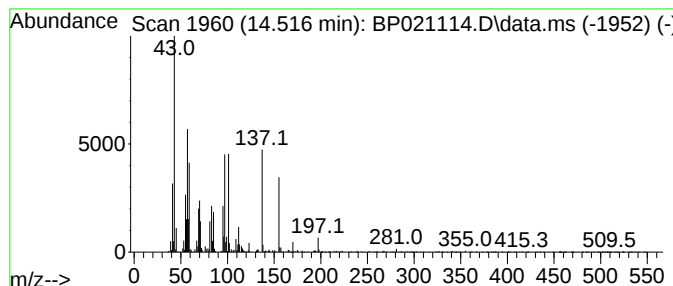
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.516	2.09 ng/ul	194030	Acenaphthene-d10	14.304

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	22
2			Ethyl octacosyl ether	438	C30H62O	1010406-37-0	10
3			Tetradecane, 1-bromo-	276	C14H29Br	000112-71-0	10
4			Succinic acid, cyclohexylmethyl ...	326	C19H34O4	1000389-54-5	9
5			2-Thiopheneacetic acid, 3-tetrad...	338	C20H34O2S	1010280-66-9	9



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072324\
Data File : BP021114.D
Acq On : 24 Jul 2024 01:51
Operator : MA/JU
Sample : P3238-06
Misc :
ALS Vial : 21 Sample Multiplier: 1

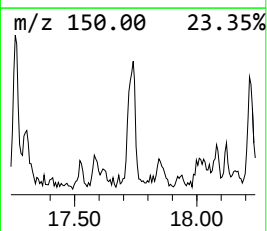
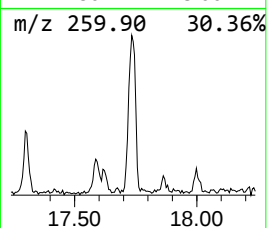
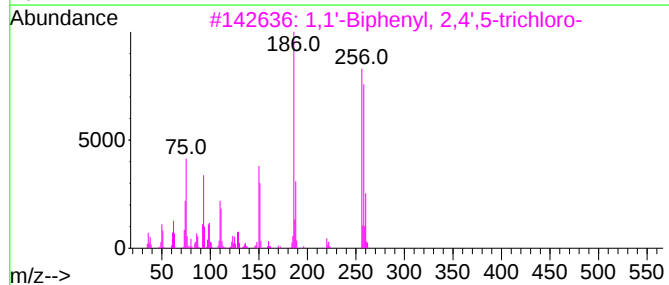
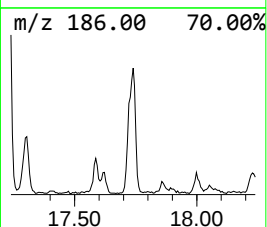
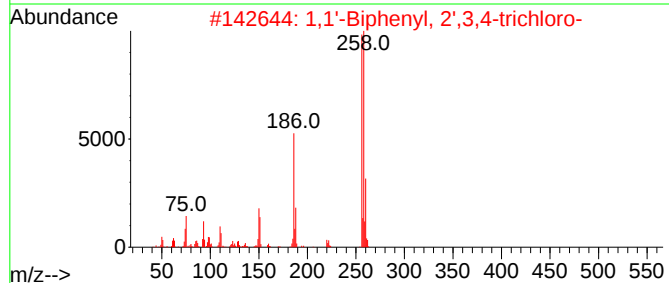
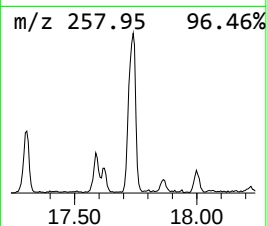
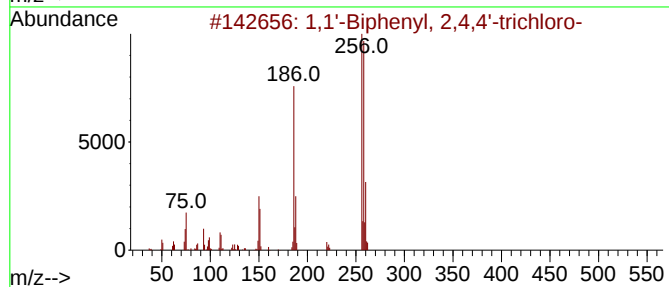
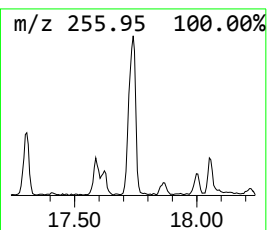
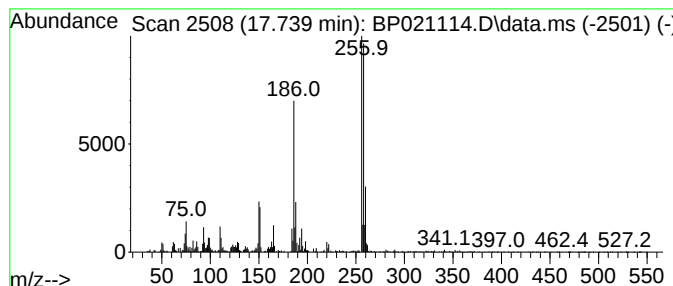
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 1,1'-Biphenyl, 2,4,4'-trich... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
17.739	2.72 ng/ul	324619	Phenanthrene-d10	17.116	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,4,4'-trichloro-	256	C12H7Cl3	007012-37-5	99
2	1,1'-Biphenyl, 2',3,4-trichloro-	256	C12H7Cl3	038444-86-9	99
3	1,1'-Biphenyl, 2,4',5-trichloro-	256	C12H7Cl3	016606-02-3	99
4	1,1'-Biphenyl, 2,3,3'-trichloro-	256	C12H7Cl3	038444-84-7	99
5	1,1'-Biphenyl, 2,3',5-trichloro-	256	C12H7Cl3	038444-81-4	99



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072324\
Data File : BP021114.D
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Sample : P3238-06
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Instrument :
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ClientSampleId :
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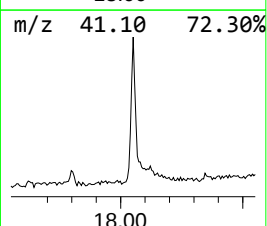
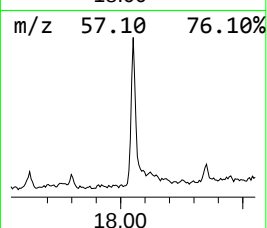
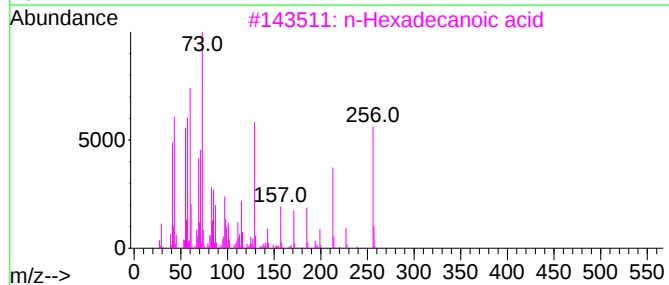
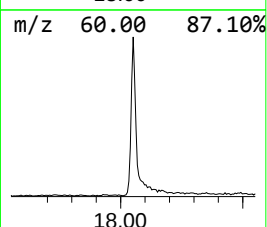
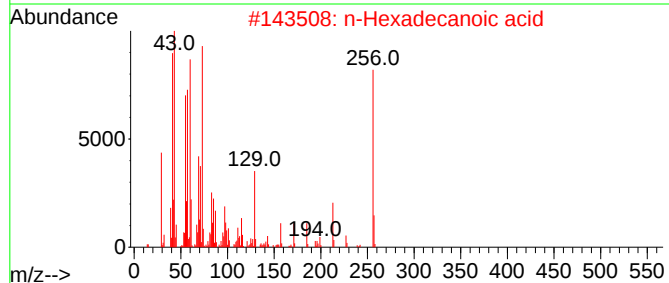
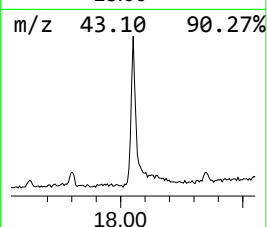
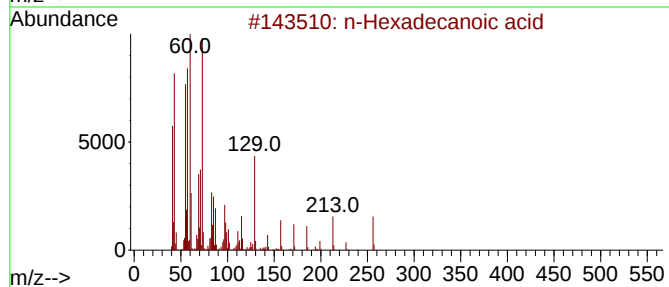
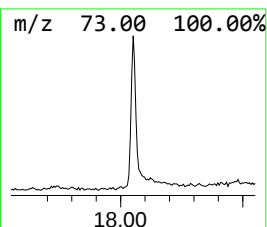
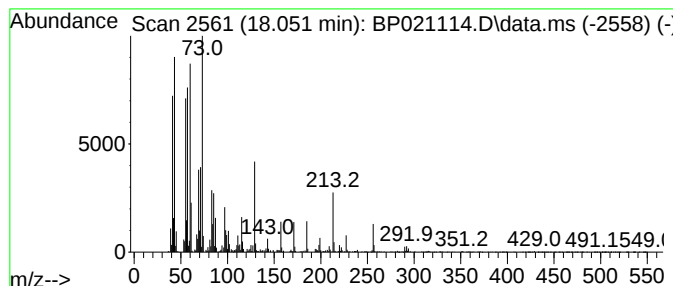
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 n-Hexadecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.051	2.78 ng/ul	331759	Phenanthrene-d10	17.116

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	97
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	95
5			Tridecanoic acid	214	C13H26O2	000638-53-9	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072324\
Data File : BP021114.D
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Sample : P3238-06
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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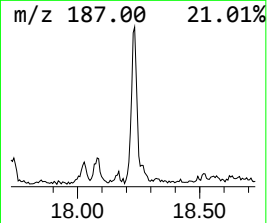
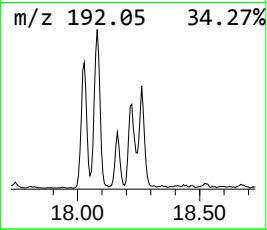
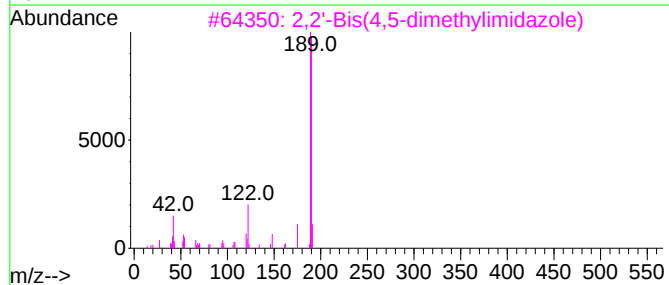
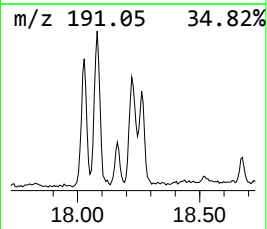
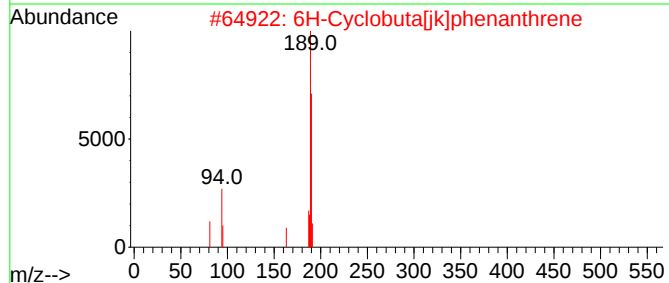
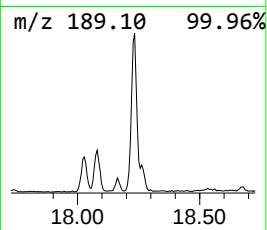
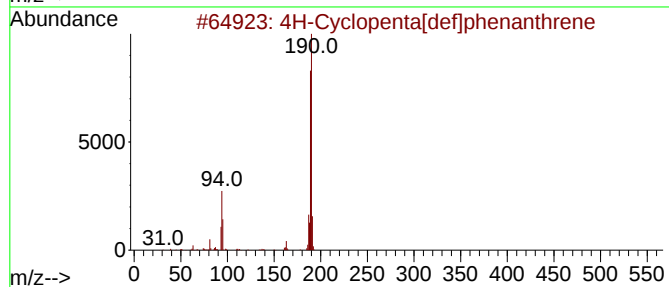
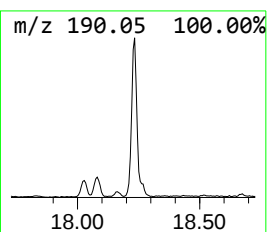
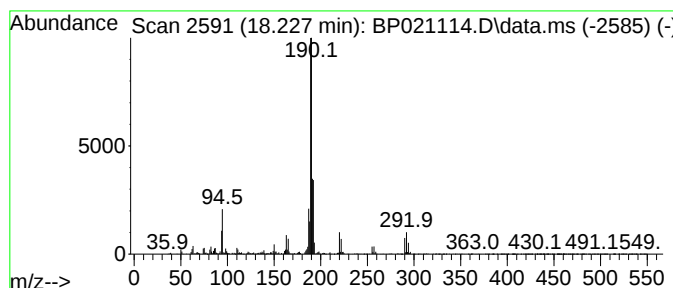
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 4H-Cyclopenta[def]phenanthrene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.227	4.91 ng/ul	586719	Phenanthrene-d10	17.116

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	81
2			6H-Cyclobuta[jk]phenanthrene	190	C15H10	083469-43-6	72
3			2,2'-Bis(4,5-dimethylimidazole)	190	C10H14N4	069286-06-2	49
4			Thiophene-3-carboxaldehyde, 5-ch...	190	C5H4BClO3S	036155-87-0	47
5			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	46



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072324\
Data File : BP021114.D
Acq On : 24 Jul 2024 01:51
Operator : MA/JU
Sample : P3238-06
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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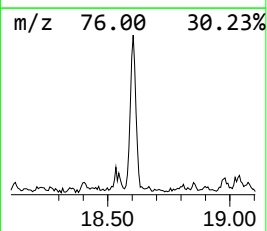
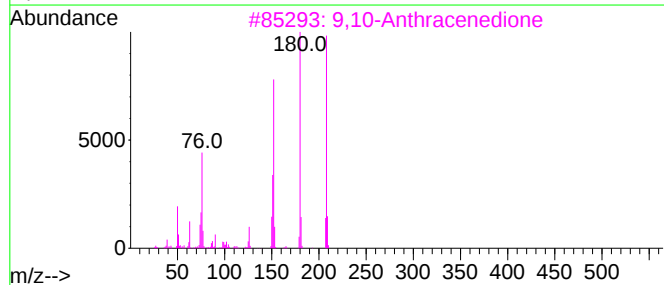
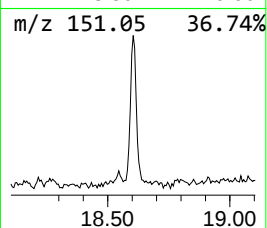
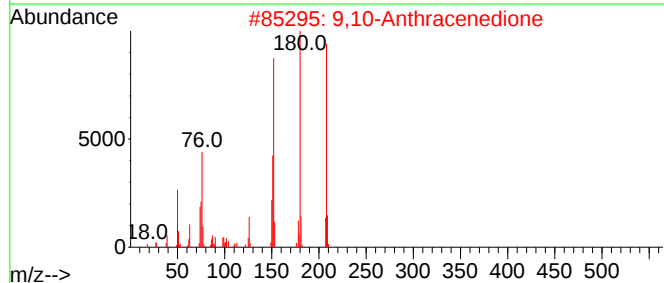
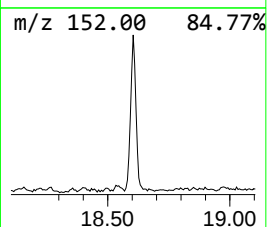
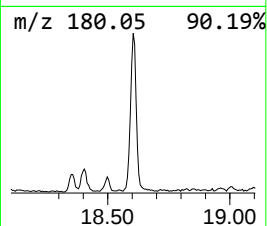
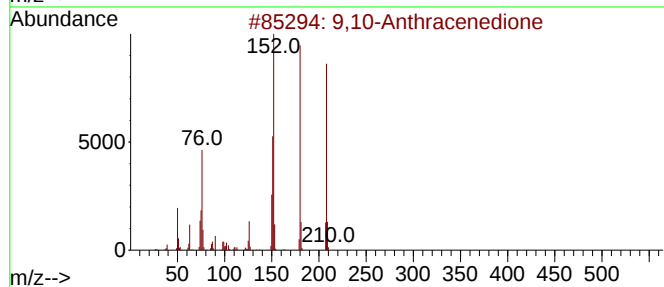
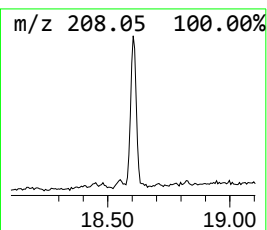
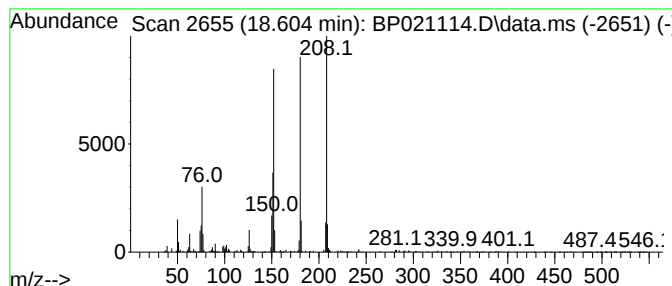
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 9,10-Anthracenedione Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.604	2.90 ng/ul	345818	Phenanthrene-d10	17.116

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	9,10-Anthracenedione			208	C14H8O2	000084-65-1	98
2	9,10-Anthracenedione			208	C14H8O2	000084-65-1	98
3	9,10-Anthracenedione			208	C14H8O2	000084-65-1	96
4	9,10-Anthracenedione			208	C14H8O2	000084-65-1	95
5	1H-Phenalen-1-one			180	C13H8O	000548-39-0	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072324\
Data File : BP021114.D
Acq On : 24 Jul 2024 01:51
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Misc :
ALS Vial : 21 Sample Multiplier: 1

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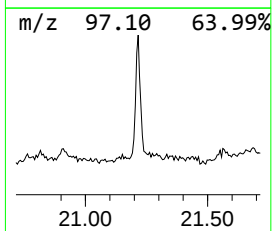
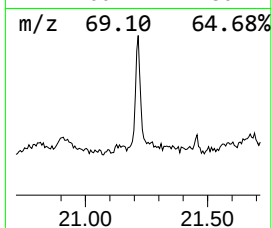
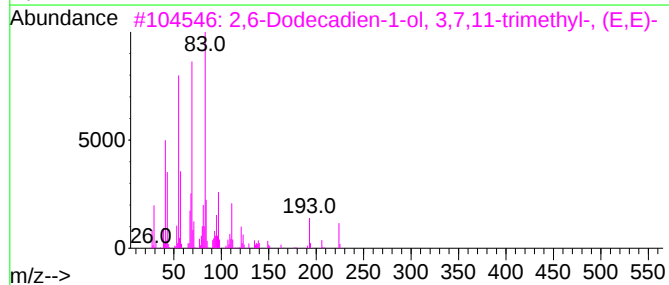
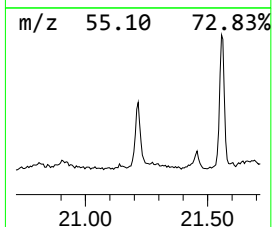
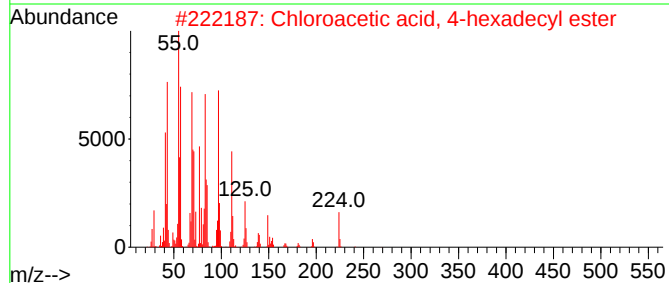
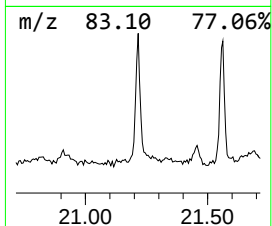
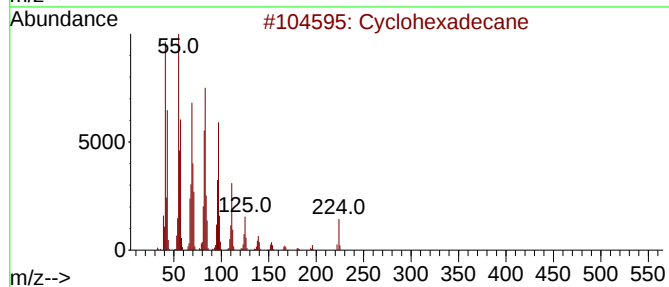
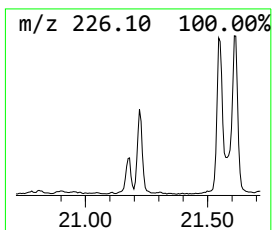
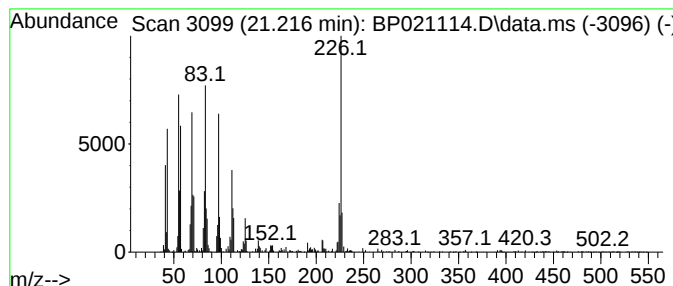
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 (DEL) Alkane: Cyclic21.215 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.215	3.16 ng/ul	912338	Chrysene-d12	21.568

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexadecane	224	C16H32	000295-65-8	58
2			Chloroacetic acid, 4-hexadecyl e...	318	C18H35ClO2	1000282-93-2	38
3			2,6-Dodecadien-1-ol, 3,7,11-trim...	224	C15H28O	020576-56-1	35
4			Pentafluoropropionic acid, 4-hex...	388	C19H33F5O2	1000283-04-1	27
5			3-Hexadecanol acetate	284	C18H36O2	051354-27-9	27



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072324\
Data File : BP021114.D
Acq On : 24 Jul 2024 01:51
Operator : MA/JU
Sample : P3238-06
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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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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1-Phenoxypropan...	11.187	4.0	ng/ul	263174	2	10.428	1318290	20.0
2-Pentanone, 4-...	14.516	2.1	ng/ul	194030	3	14.304	1855400	20.0
1,1'-Biphenyl, ...	17.739	2.7	ng/ul	324619	4	17.116	2387650	20.0
n-Hexadecanoic ...	18.051	2.8	ng/ul	331759	4	17.116	2387650	20.0
4H-Cyclopenta[d...]	18.227	4.9	ng/ul	586719	4	17.116	2387650	20.0
9,10-Anthracene...	18.604	2.9	ng/ul	345818	4	17.116	2387650	20.0
(DEL) Alkane: C...	21.215	3.2	ng/ul	912338	5	21.568	5772490	20.0