

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071524\
 Data File : BP020959.D
 Acq On : 15 Jul 2024 15:06
 Operator : MA/JU
 Sample : P3177-21
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 A4DN2

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: BP020959.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.234	39	42	46	rBV	23742	27241	0.12%	0.010%
2	6.064	518	523	527	rBV2	18641	30349	0.14%	0.011%
3	6.875	649	661	662	rBV	497052	797518	3.59%	0.291%
4	7.011	678	684	690	rBV	403185	600064	2.70%	0.219%
5	7.099	693	699	712	rVB	2002169	3106265	13.99%	1.134%
6	7.216	712	719	720	rBV	546779	776520	3.50%	0.283%
7	7.246	720	724	733	rVB	1931637	3077987	13.86%	1.123%
8	7.669	790	796	797	rBV	669903	832840	3.75%	0.304%
9	7.699	797	801	814	rVB	3517015	4989086	22.47%	1.821%
10	8.011	847	854	872	rVB	2877198	4771674	21.49%	1.741%
11	8.387	911	918	925	rBV	572176	1021805	4.60%	0.373%
12	8.463	925	931	942	rVB	2693563	4009690	18.06%	1.463%
13	8.816	983	991	1001	rBV	455809	787706	3.55%	0.287%
14	9.205	1054	1057	1062	rVB2	19212	29428	0.13%	0.011%
15	9.387	1074	1088	1102	rBV	3243389	6220932	28.02%	2.270%
16	9.563	1106	1118	1130	rVV2	1139644	2508572	11.30%	0.916%
17	9.675	1131	1137	1147	rVB	119682	231302	1.04%	0.084%
18	10.093	1196	1208	1223	rBV2	4006776	8066464	36.33%	2.944%
19	10.281	1238	1240	1247	rVB5	13426	22917	0.10%	0.008%
20	10.428	1258	1265	1268	rVV	774788	1333446	6.01%	0.487%
21	10.481	1268	1274	1284	rVV	4609000	7814867	35.20%	2.852%
22	10.581	1284	1291	1309	rVV	475472	939063	4.23%	0.343%
23	10.710	1309	1313	1318	rVB3	16944	27016	0.12%	0.010%
24	11.175	1384	1392	1399	rBV	62267	134163	0.60%	0.049%
25	11.746	1480	1489	1499	rBV	615183	1159317	5.22%	0.423%
26	11.828	1499	1503	1512	rVB8	11412	33110	0.15%	0.012%
27	11.981	1525	1529	1534	rBV2	14622	28869	0.13%	0.011%
28	12.104	1541	1550	1556	rBV	5781849	9268951	41.75%	3.383%
29	12.210	1562	1568	1575	rVB2	23406	50445	0.23%	0.018%
30	12.304	1578	1584	1592	rVB	59809	95565	0.43%	0.035%
31	12.716	1646	1654	1662	rBV	982057	1519344	6.84%	0.554%
32	12.804	1662	1669	1679	rVB	1425401	2640876	11.90%	0.964%
33	13.087	1712	1717	1721	rVB	19542	26659	0.12%	0.010%
34	13.228	1737	1741	1749	rBV2	23970	45602	0.21%	0.017%
35	13.334	1756	1759	1769	rVB3	25847	47584	0.21%	0.017%
36	13.710	1813	1823	1831	rBV	1145005	1685584	7.59%	0.615%

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Peak Location: TOP

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

37	13.893	1845	1854	1861	rBV	5717802	9277086	41.79%	3.386%
38	14.028	1865	1877	1888	rVB2	4503836	8519487	38.37%	3.109%
39	14.175	1896	1902	1908	rVB6	16033	31326	0.14%	0.011%
40	14.304	1914	1924	1930	rBV	1076214	1893739	8.53%	0.691%

41	14.375	1930	1936	1941	rVV	7230199	10674874	48.08%	3.896%
42	14.422	1941	1944	1949	rVB	101943	127235	0.57%	0.046%
43	14.522	1951	1961	1965	rBV	2218212	5504112	24.79%	2.009%
44	14.569	1965	1969	1978	rVB3	789638	1539834	6.94%	0.562%
45	14.710	1985	1993	2005	rVB2	4402342	7296889	32.87%	2.663%

46	14.946	2028	2033	2043	rVB	56076	97010	0.44%	0.035%
47	15.145	2055	2067	2074	rBV	6246627	10007055	45.08%	3.652%
48	15.316	2089	2096	2099	rBV	1651942	2650872	11.94%	0.967%
49	15.369	2099	2105	2116	rVV2	6227016	13259704	59.73%	4.839%
50	15.463	2116	2121	2131	rVV	508412	743727	3.35%	0.271%

51	15.551	2134	2136	2143	rVB7	13590	27323	0.12%	0.010%
52	15.628	2144	2149	2154	rBV2	87543	121887	0.55%	0.044%
53	15.881	2187	2192	2196	rBV6	28477	49733	0.22%	0.018%
54	15.928	2196	2200	2204	rVV3	28181	46739	0.21%	0.017%
55	15.969	2204	2207	2211	rVB3	39991	51995	0.23%	0.019%

56	16.063	2216	2223	2236	rVB4	53897	139635	0.63%	0.051%
57	16.281	2249	2260	2268	rBV	2945858	4305924	19.40%	1.571%
58	16.640	2318	2321	2327	rVB2	18943	25609	0.12%	0.009%
59	16.763	2335	2342	2355	rBV	3662662	5655429	25.47%	2.064%
60	16.910	2362	2367	2378	rVB4	97607	218074	0.98%	0.080%

61	17.051	2384	2391	2394	rBV7	18424	41170	0.19%	0.015%
62	17.110	2394	2401	2404	rVV	1349687	2234378	10.06%	0.815%
63	17.151	2404	2408	2415	rVV	2491450	4632581	20.87%	1.691%
64	17.222	2415	2420	2422	rVV2	1841359	2654767	11.96%	0.969%
65	17.257	2422	2426	2439	rVB	7347765	10614988	47.81%	3.874%

66	17.410	2447	2452	2457	rVB2	26664	52562	0.24%	0.019%
67	17.534	2467	2473	2484	rVB4	42426	114232	0.51%	0.042%
68	17.745	2505	2509	2515	rBV2	119008	164634	0.74%	0.060%
69	17.810	2516	2520	2527	rVB	287501	327926	1.48%	0.120%
70	18.051	2556	2561	2566	rBV	504258	793159	3.57%	0.289%

71	18.116	2566	2572	2587	rVB	5591883	9205914	41.47%	3.360%
72	18.434	2621	2626	2638	rBV4	97758	230641	1.04%	0.084%
73	18.528	2638	2642	2647	rVV	86610	126585	0.57%	0.046%
74	18.598	2649	2654	2665	rVV	584758	897587	4.04%	0.328%
75	18.834	2690	2694	2698	rVB2	31496	45674	0.21%	0.017%

76	18.951	2710	2714	2717	rBV5	25944	45717	0.21%	0.017%
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77	18.998	2719	2722	2724	rVV3	47300	55168	0.25%	0.020%
78	19.028	2724	2727	2731	rVB	201270	239593	1.08%	0.087%
79	19.175	2749	2752	2756	rVB	98705	109489	0.49%	0.040%
80	19.257	2759	2766	2778	rVB	5999606	9144413	41.19%	3.337%
81	19.386	2783	2788	2796	rBV2	215410	444160	2.00%	0.162%
82	19.598	2819	2824	2833	rBV	2308059	3407273	15.35%	1.243%
83	19.722	2842	2845	2849	rBV3	60541	76614	0.35%	0.028%
84	20.328	2945	2948	2951	rBV	158728	165703	0.75%	0.060%
85	20.581	2987	2991	2996	rBV	1447904	1705530	7.68%	0.622%
86	21.228	3097	3101	3110	rVB	360906	615377	2.77%	0.225%
87	21.386	3125	3128	3134	rVB	282910	327128	1.47%	0.119%
88	21.475	3136	3143	3148	rBV	15435731	22200741	100.00%	8.102%
89	21.551	3149	3156	3165	rVV2	4466442	9638019	43.41%	3.517%
90	22.733	3349	3357	3366	rVB	7153897	13448845	60.58%	4.908%
91	23.822	3534	3542	3547	rBV	2680956	5838853	26.30%	2.131%
92	23.886	3547	3553	3560	rVB	2150126	4520062	20.36%	1.650%
93	24.633	3673	3680	3687	rBV	1235161	3382069	15.23%	1.234%
94	24.710	3687	3693	3706	rVB	1661857	4131385	18.61%	1.508%
95	24.869	3711	3720	3729	rBV	955275	2564057	11.55%	0.936%
96	28.674	4359	4367	4368	rBV	547399	1105663	4.98%	0.404%
97	29.780	4545	4555	4556	rBV	1310210	2770150	12.48%	1.011%
98	29.809	4556	4560	4577	rVB	2018257	4912901	22.13%	1.793%

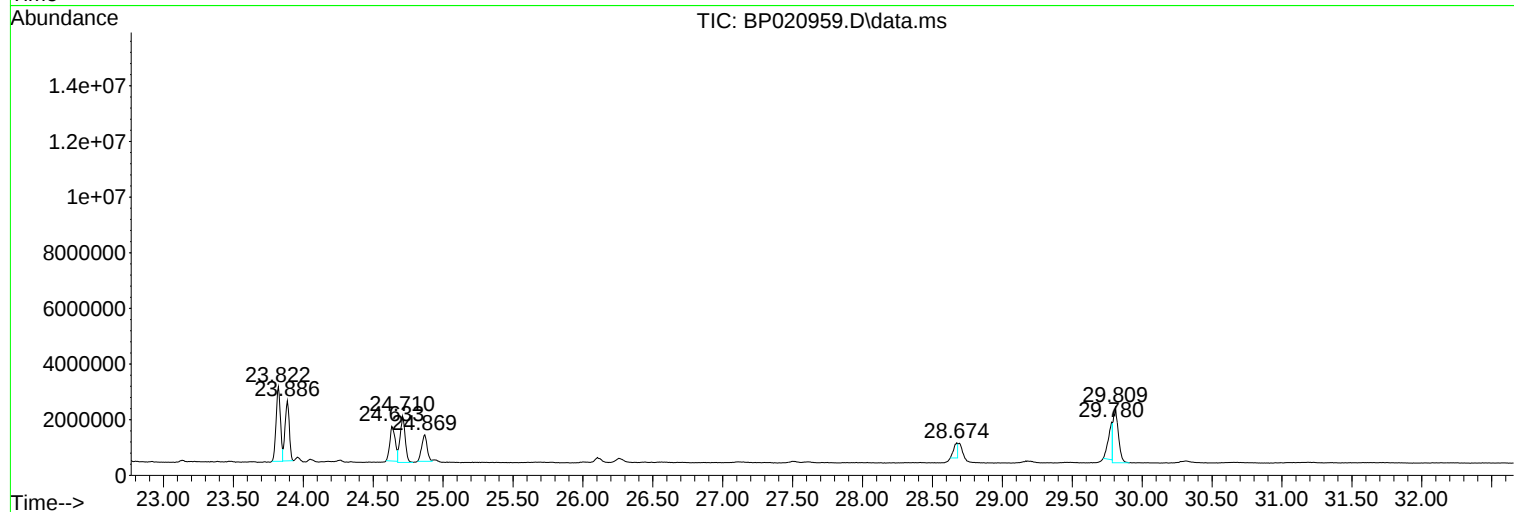
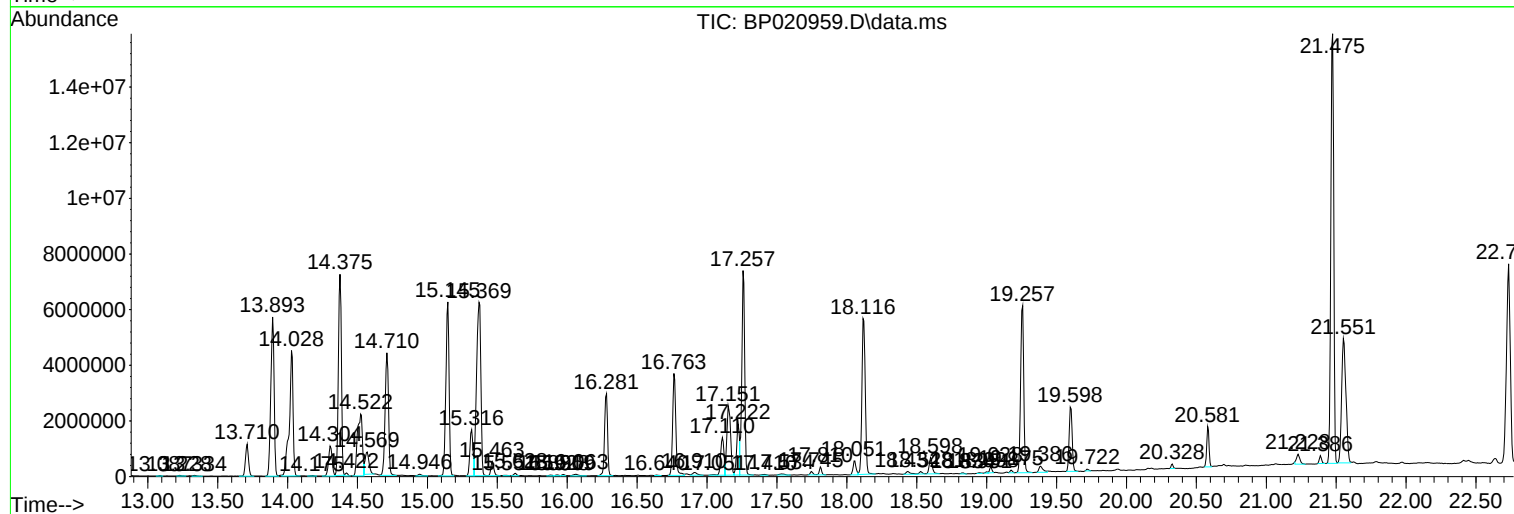
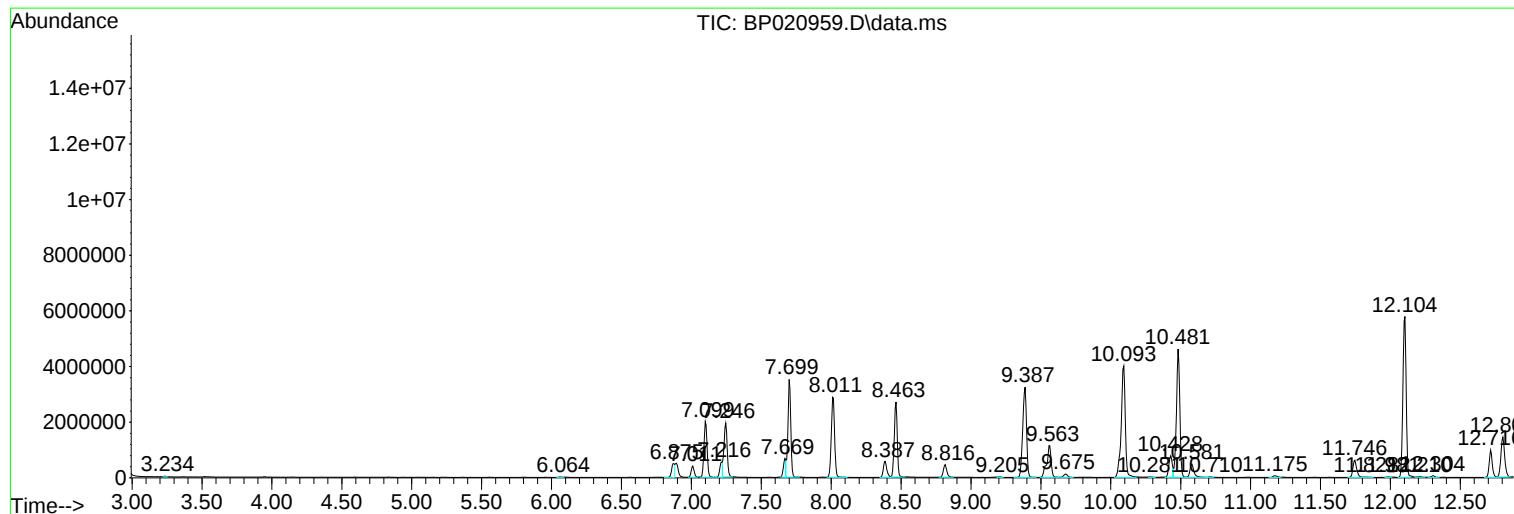
Sum of corrected areas: 274009827

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Instrument :
BNA_P
ClientSampleId :
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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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Sample : P3177-21
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Instrument :
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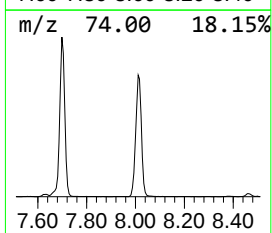
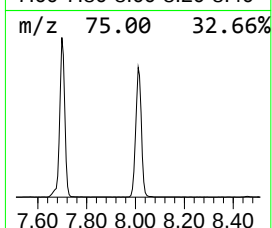
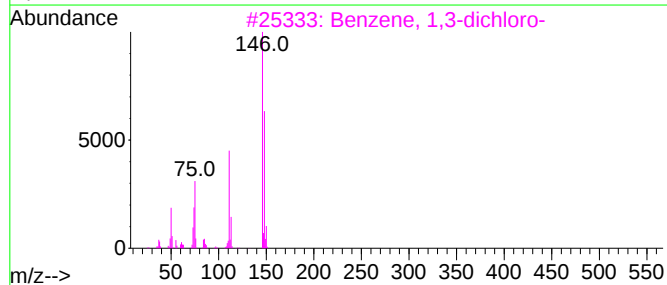
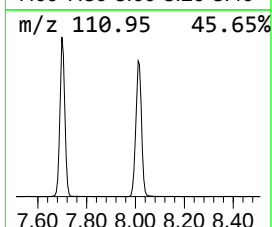
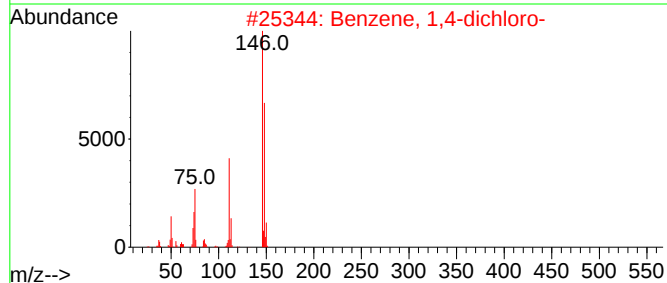
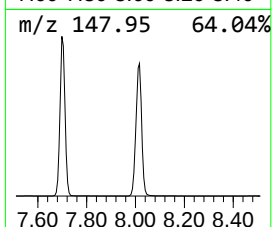
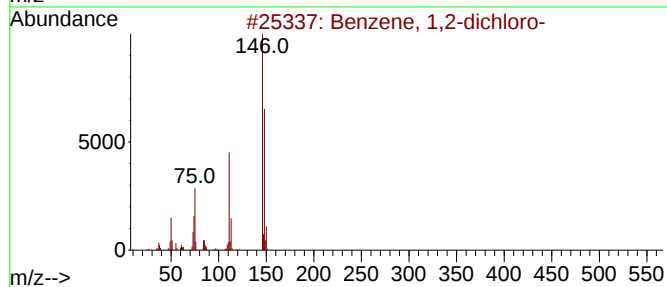
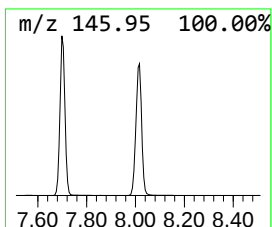
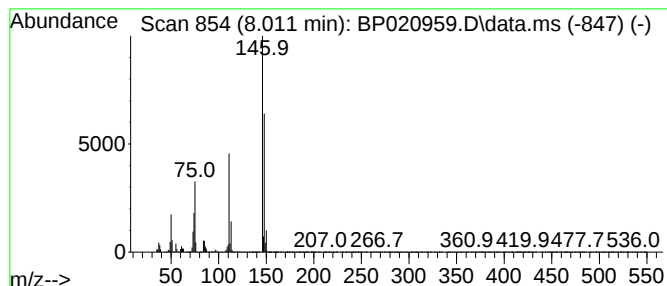
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Benzene, 1,2-dichloro- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.011	114.59 ng/ul	4771670	1,4-Dichlorobenzene-d4	7.669

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	98
2			Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	98
3			Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	98
4			Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	97
5			Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	97



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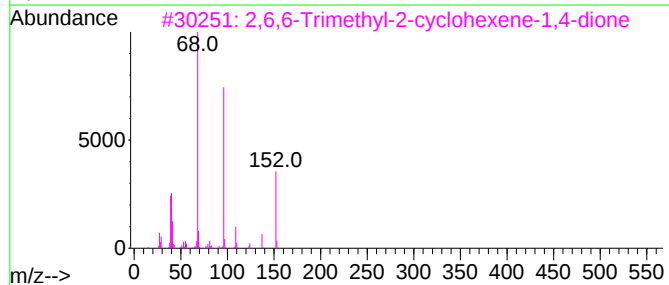
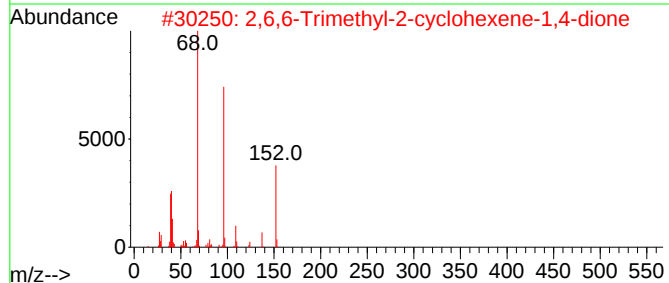
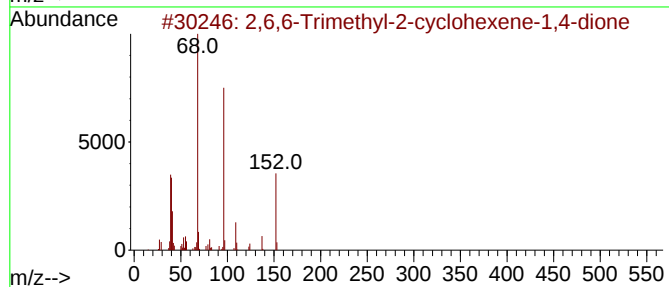
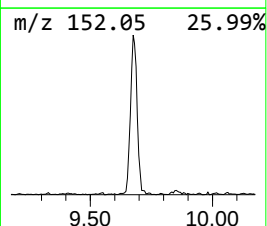
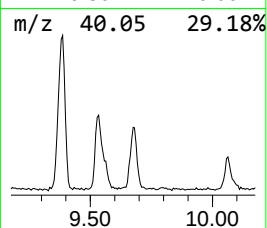
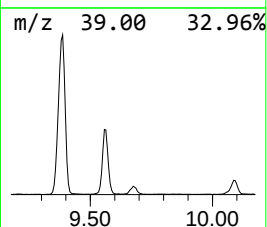
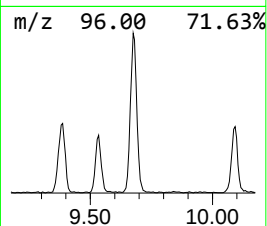
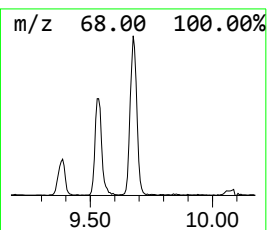
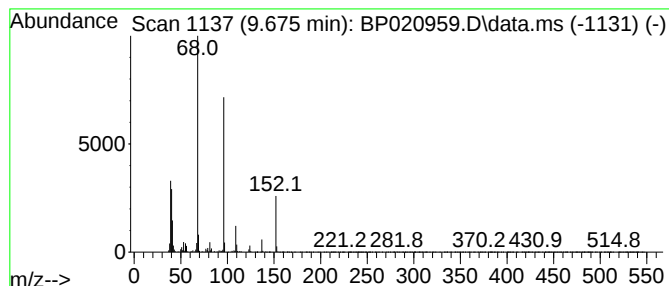
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 2,6,6-Trimethyl-2-cyclohexe... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.675	3.47 ng/ul	231302	Naphthalene-d8	10.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,6,6-Trimethyl-2-cyclohexene-1,...	152	C9H12O2	001125-21-9	96		
2	2,6,6-Trimethyl-2-cyclohexene-1,...	152	C9H12O2	001125-21-9	94		
3	2,6,6-Trimethyl-2-cyclohexene-1,...	152	C9H12O2	001125-21-9	94		
4	2,6,6-Trimethyl-2-cyclohexene-1,...	152	C9H12O2	001125-21-9	90		
5	2-Cyclohexen-1-one	96	C6H8O	000930-68-7	53		



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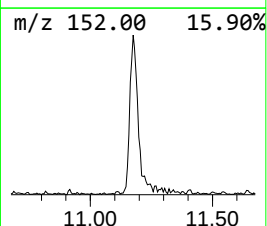
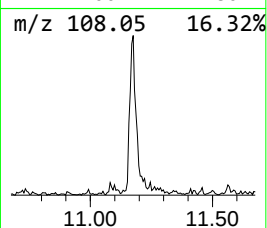
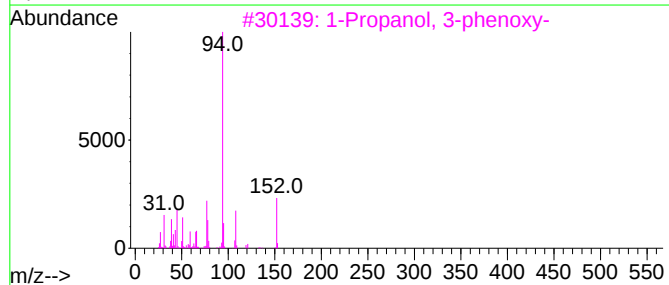
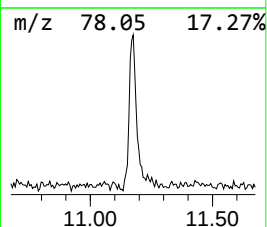
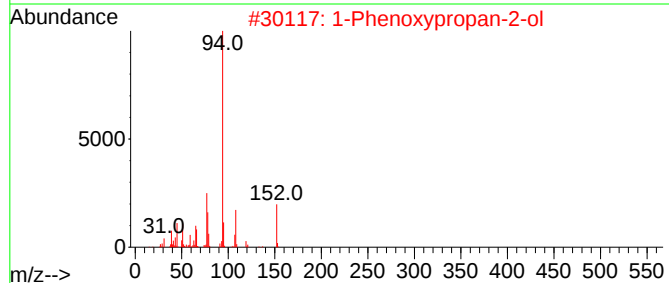
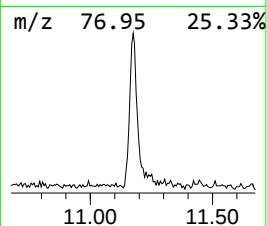
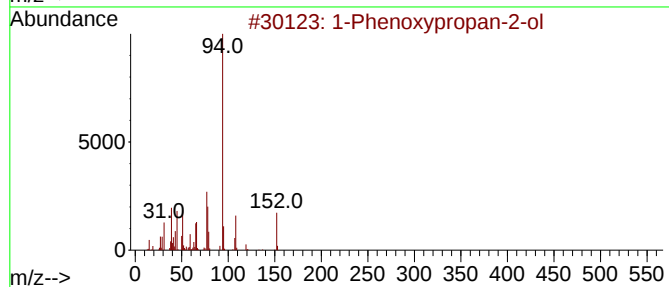
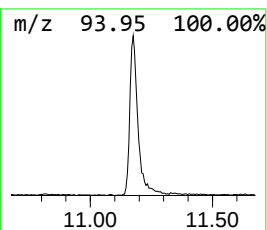
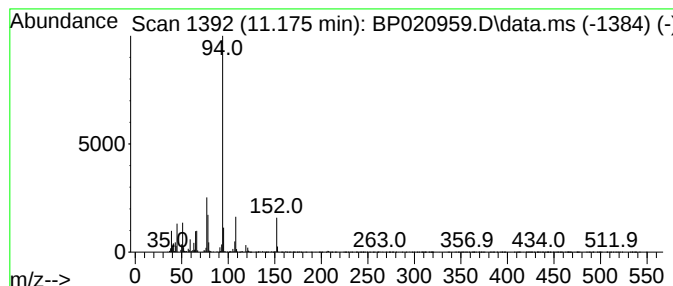
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 1-Phenoxypropan-2-ol Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.175	2.01 ng/ul	134163	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-Propanol, 3-phenoxy-	152	C9H12O2	006180-61-6	93
4			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	87
5			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071524\
Data File : BP020959.D
Acq On : 15 Jul 2024 15:06
Operator : MA/JU
Sample : P3177-21
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4DN2

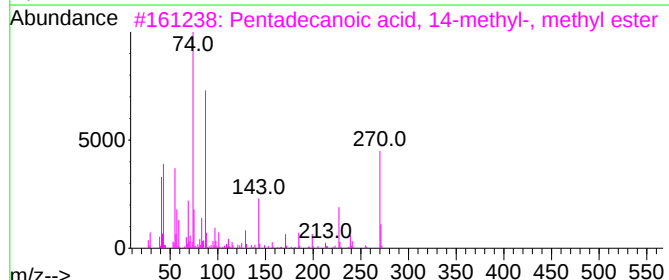
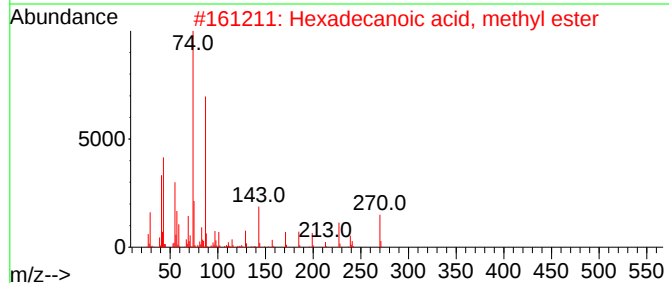
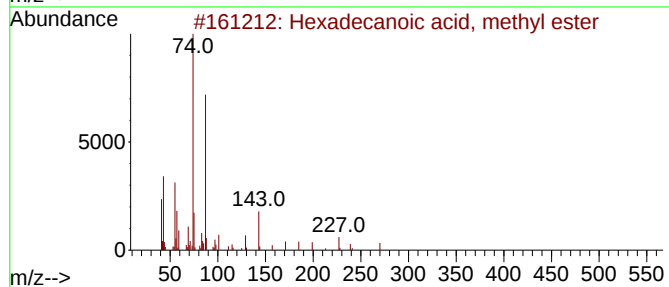
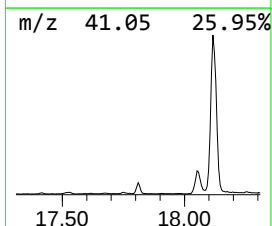
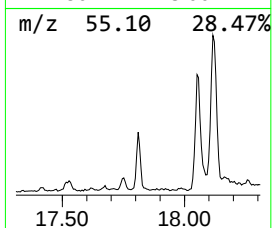
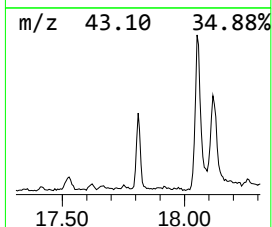
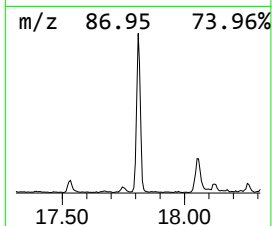
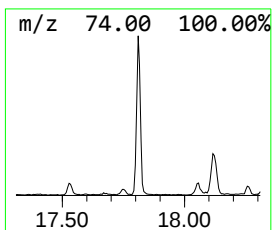
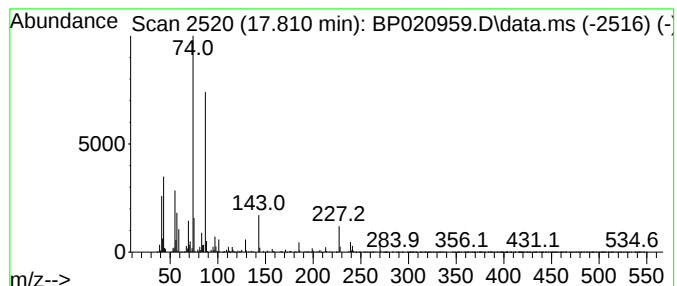
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Hexadecanoic acid, methyl e... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.810	2.94 ng/ul	327926	Phenanthrene-d10	17.110

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	94
2			Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	94
3			Pentadecanoic acid, 14-methyl-, ...	270	C17H34O2	005129-60-2	94
4			Pentadecanoic acid, 14-methyl-, ...	270	C17H34O2	005129-60-2	93
5			Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071524\
Data File : BP020959.D
Acq On : 15 Jul 2024 15:06
Operator : MA/JU
Sample : P3177-21
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4DN2

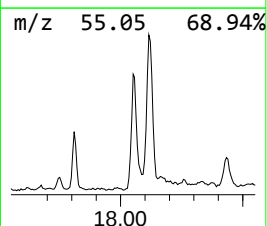
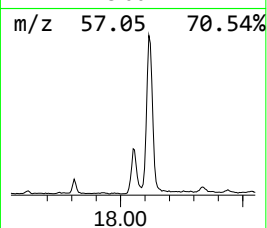
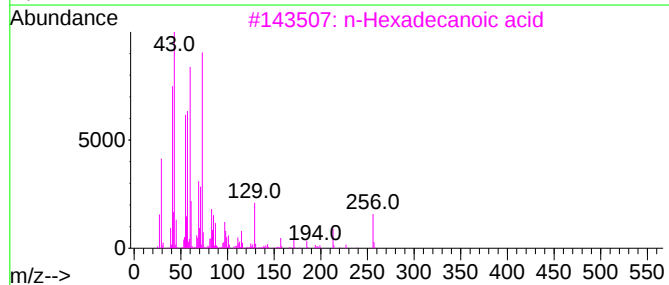
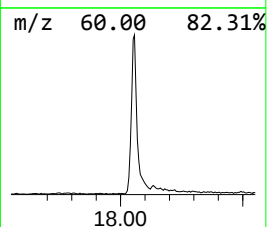
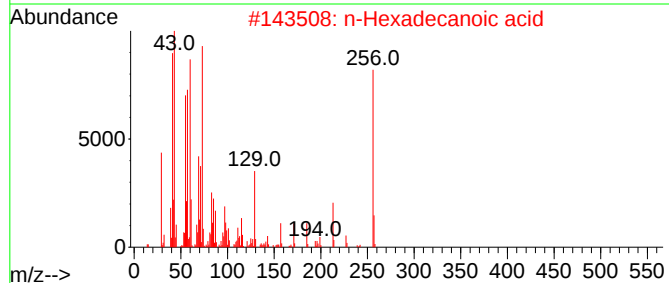
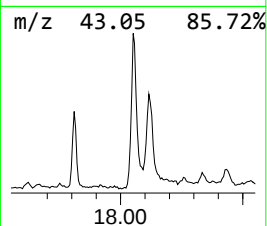
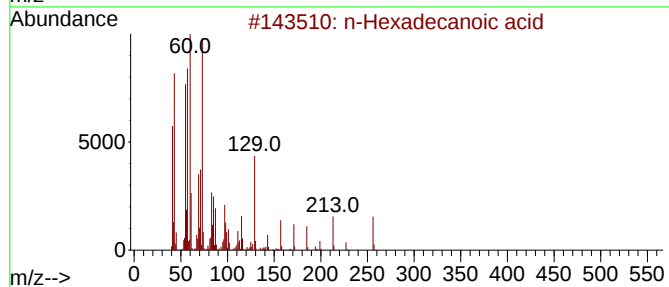
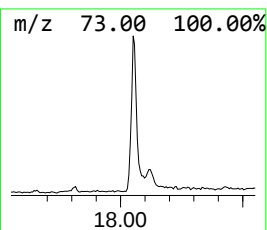
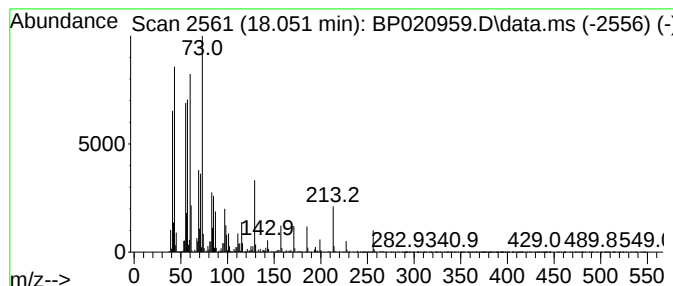
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.051	7.10 ng/ul	793159	Phenanthrene-d10	17.110

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			Tetradecanoic acid	228	C14H28O2	000544-63-8	94
5			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071524\
Data File : BP020959.D
Acq On : 15 Jul 2024 15:06
Operator : MA/JU
Sample : P3177-21
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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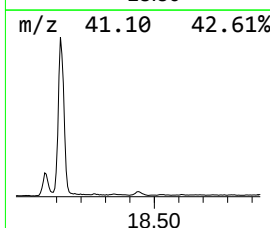
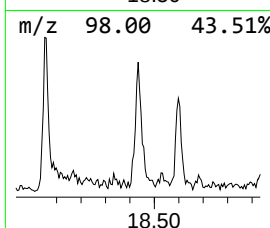
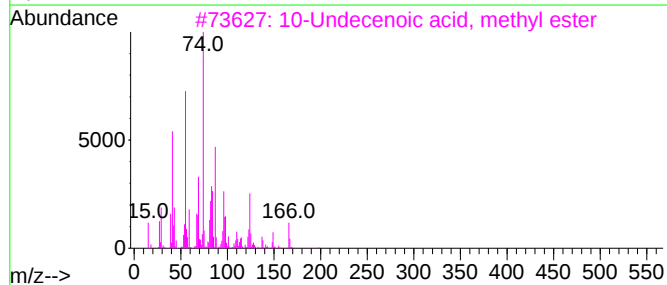
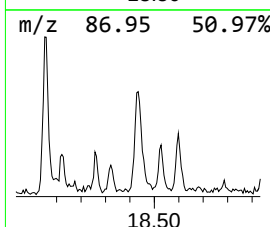
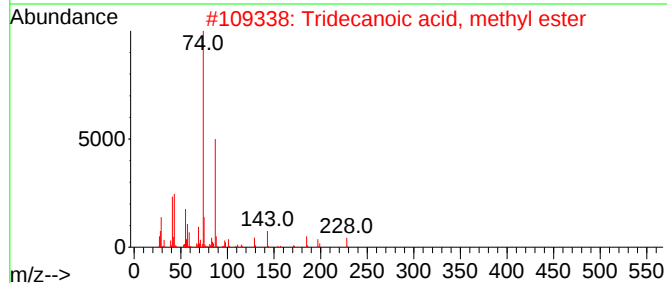
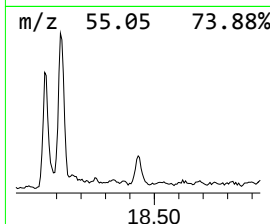
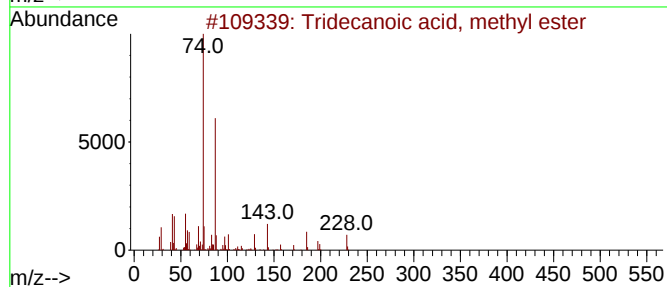
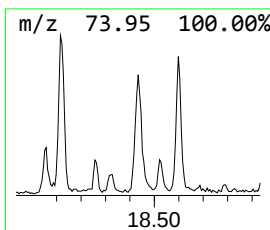
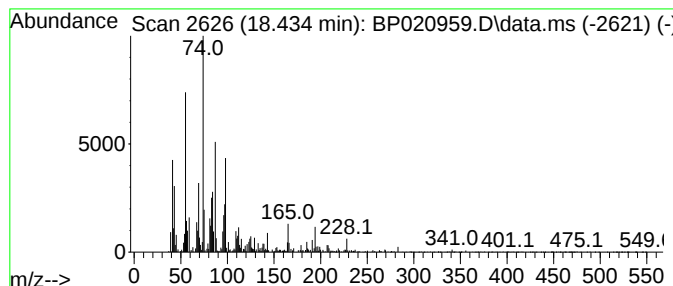
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Tridecanoic acid, methyl ester Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.434	2.06 ng/ul	230641	Phenanthrene-d10	17.110

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tridecanoic acid, methyl ester	228	C14H28O2	001731-88-0	50
2			Tridecanoic acid, methyl ester	228	C14H28O2	001731-88-0	50
3			10-Undecenoic acid, methyl ester	198	C12H22O2	000111-81-9	49
4			Tridecanoic acid, methyl ester	228	C14H28O2	001731-88-0	46
5			Dodecanoic acid, 10-methyl-, met...	228	C14H28O2	005129-65-7	45



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071524\
Data File : BP020959.D
Acq On : 15 Jul 2024 15:06
Operator : MA/JU
Sample : P3177-21
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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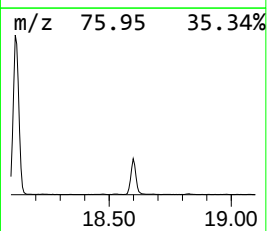
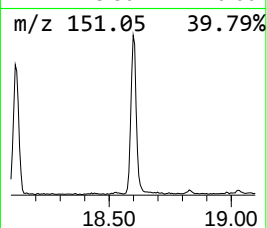
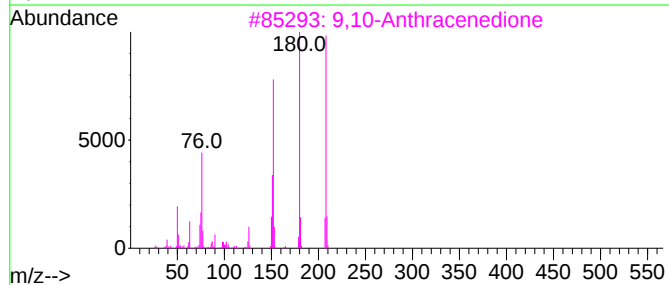
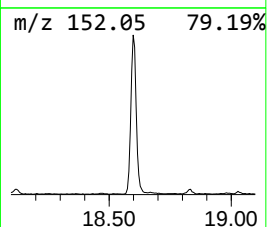
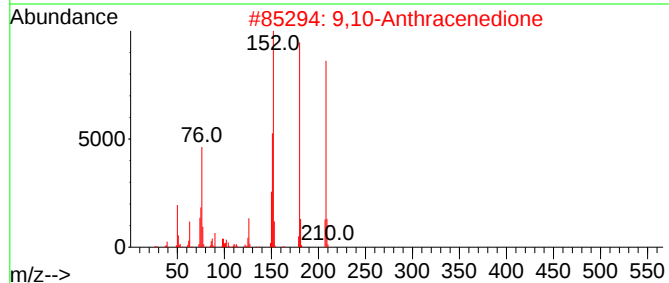
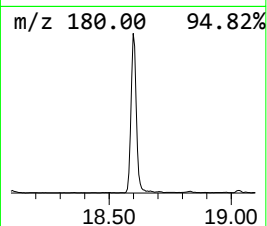
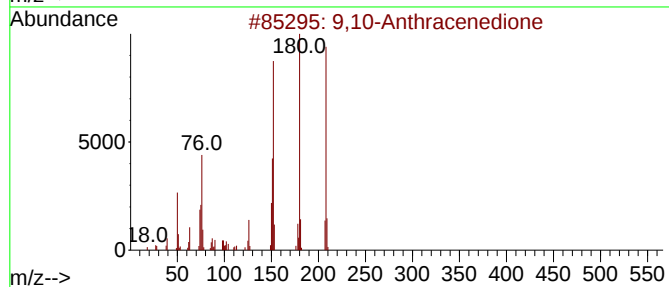
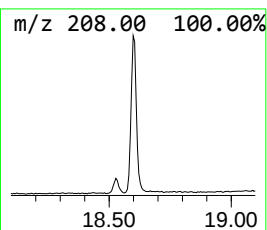
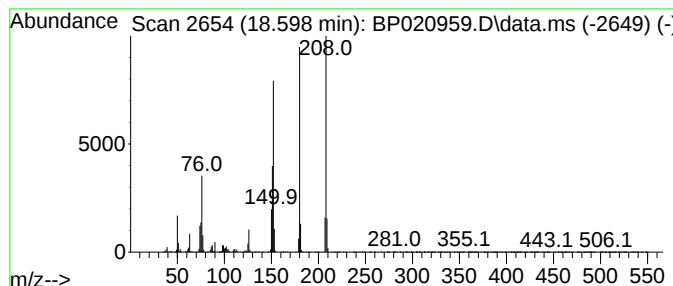
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 9,10-Anthracenedione Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.598	8.03 ng/ul	897587	Phenanthrene-d10	17.110

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	9,10-Anthracenedione			208	C14H8O2	000084-65-1	99
2	9,10-Anthracenedione			208	C14H8O2	000084-65-1	98
3	9,10-Anthracenedione			208	C14H8O2	000084-65-1	98
4	9,10-Anthracenedione			208	C14H8O2	000084-65-1	94
5	9H-Fluoren-9-one			180	C13H8O	000486-25-9	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071524\
Data File : BP020959.D
Acq On : 15 Jul 2024 15:06
Operator : MA/JU
Sample : P3177-21
Misc :
ALS Vial : 6 Sample Multiplier: 1

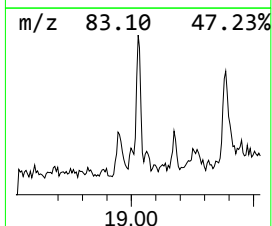
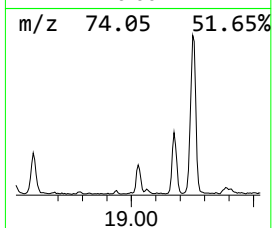
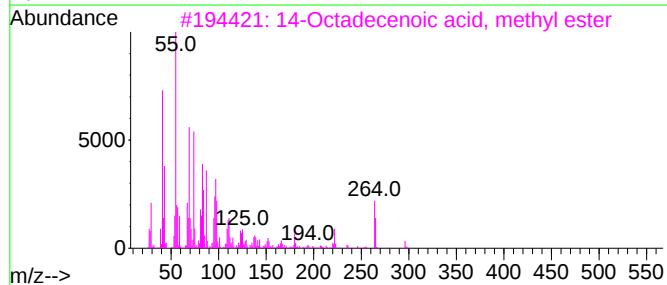
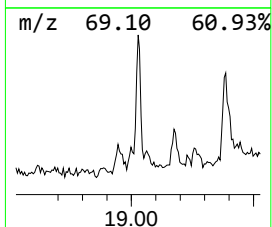
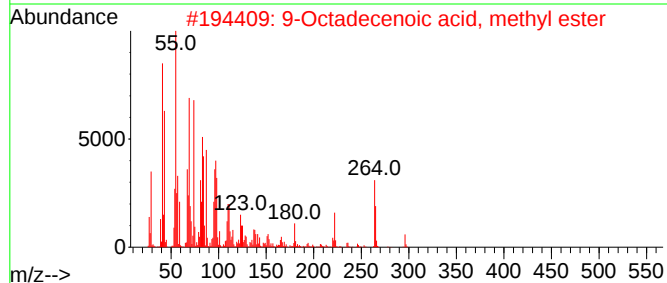
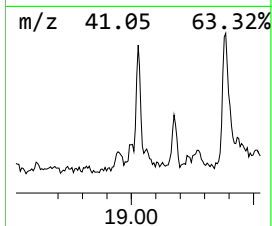
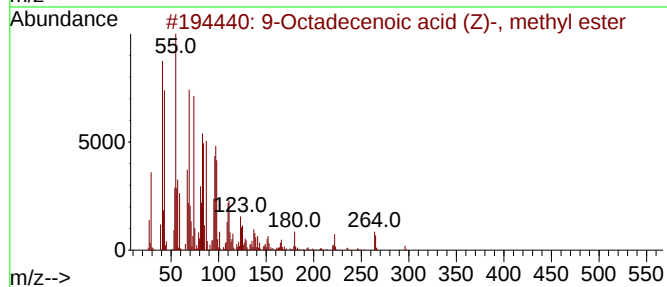
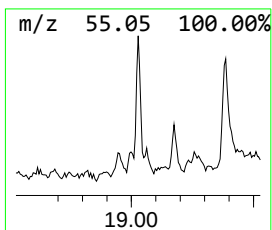
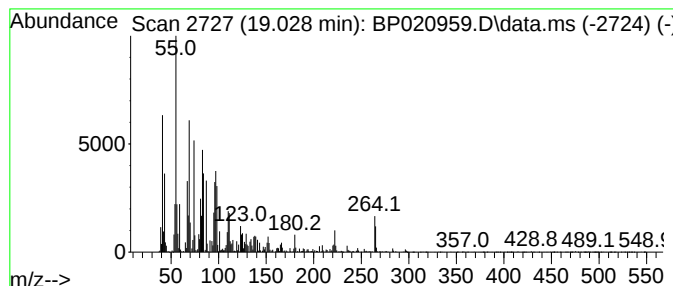
Instrument :
BNA_P
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 8 9-Octadecenoic acid (Z)-, m... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD		R.T.	
19.028	2.14 ng/ul	239593	Phenanthrene-d10		17.110	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	9-Octadecenoic acid (Z)-, methyl...		296	C19H36O2	000112-62-9	99
2	9-Octadecenoic acid, methyl ester		296	C19H36O2	002462-84-2	99
3	14-Octadecenoic acid, methyl ester		296	C19H36O2	056554-48-4	97
4	11-Octadecenoic acid, methyl ester		296	C19H36O2	052380-33-3	97
5	(Z)-15-Octadecenoic acid methyl ...		296	C19H36O2	1000427-69-3	96



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071524\
Data File : BP020959.D
Acq On : 15 Jul 2024 15:06
Operator : MA/JU
Sample : P3177-21
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4DN2

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.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
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2,6,6-Trimethyl...	9.675	3.5	ng/ul	231302	2	10.428	1333450	20.0
1-Phenoxypropan...	11.175	2.0	ng/ul	134163	2	10.428	1333450	20.0
Hexadecanoic ac...	17.810	2.9	ng/ul	327926	4	17.110	2234380	20.0
n-Hexadecanoic ...	18.051	7.1	ng/ul	793159	4	17.110	2234380	20.0
Tridecanoic aci...	18.434	2.1	ng/ul	230641	4	17.110	2234380	20.0
9,10-Anthracene...	18.598	8.0	ng/ul	897587	4	17.110	2234380	20.0
9-Octadecenoic ...	19.028	2.1	ng/ul	239593	4	17.110	2234380	20.0