

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070124\
 Data File : BP020743.D
 Acq On : 02 Jul 2024 09:26
 Operator : MA/JU
 Sample : P2963-19
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 A4DK6

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-BP062524.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP020743.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.511	253	259	270	rVB3	45440	91332	0.32%	0.019%
2	6.922	661	669	671	rBV	974419	1623831	5.66%	0.332%
3	6.946	671	673	687	rVB	1353980	2086359	7.27%	0.426%
4	7.087	690	697	703	rBV	712730	1129195	3.93%	0.231%
5	7.181	703	713	722	rVB	5743153	8473426	29.51%	1.730%
6	7.317	731	736	747	rVB	3308930	5680973	19.79%	1.160%
7	7.781	809	815	824	rVB	7965932	13064496	45.51%	2.668%
8	8.093	858	868	878	rBV	7929350	13090159	45.59%	2.673%
9	8.463	923	931	938	rBV	1134746	1900527	6.62%	0.388%
10	8.558	938	947	969	rVB	5764274	10275953	35.79%	2.098%
11	8.811	982	990	997	rVB	4409026	7394873	25.76%	1.510%
12	8.893	997	1004	1007	rBV	950466	1482128	5.16%	0.303%
13	8.934	1007	1011	1019	rVB	1529087	2287651	7.97%	0.467%
14	9.275	1064	1069	1077	rVB3	33866	67960	0.24%	0.014%
15	9.469	1087	1102	1117	rBV	6651430	13168312	45.87%	2.689%
16	9.610	1117	1126	1127	rVV	827937	1203257	4.19%	0.246%
17	9.634	1127	1130	1141	rVV	1596743	2446903	8.52%	0.500%
18	9.758	1145	1151	1158	rVV	291670	525015	1.83%	0.107%
19	9.934	1173	1181	1194	rBV	6350417	10494449	36.55%	2.143%
20	10.175	1209	1222	1235	rBV2	4394242	10101292	35.18%	2.063%
21	10.510	1261	1279	1286	rBV	1126531	2012704	7.01%	0.411%
22	10.658	1296	1304	1319	rBV	824491	1440627	5.02%	0.294%
23	10.834	1322	1334	1342	rBV	7213381	11932055	41.56%	2.436%
24	11.257	1397	1406	1425	rBV2	118193	303956	1.06%	0.062%
25	11.799	1491	1498	1514	rVB	2254360	3609701	12.57%	0.737%
26	12.075	1539	1545	1549	rBV	79745	141875	0.49%	0.029%
27	12.169	1554	1561	1567	rVV	5009149	6935719	24.16%	1.416%
28	12.216	1567	1569	1576	rVV3	88325	147494	0.51%	0.030%
29	12.593	1626	1633	1645	rVB	135610	198619	0.69%	0.041%
30	12.793	1659	1667	1673	rVV	9833713	15134078	52.71%	3.090%
31	12.863	1673	1679	1693	rVB	2088387	3653784	12.73%	0.746%
32	13.181	1727	1733	1737	rBV2	41605	62272	0.22%	0.013%
33	13.287	1746	1751	1760	rBV	52559	90084	0.31%	0.018%
34	13.369	1760	1765	1770	rVV	39321	65466	0.23%	0.013%
35	13.416	1770	1773	1782	rVB2	30553	52419	0.18%	0.011%
36	13.793	1829	1837	1850	rBV	1918893	3164038	11.02%	0.646%

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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

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062524.MA.M
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37	13.975	1859	1868	1872	rBV	9549846	15004536	52.26%	3.064%
38	14.063	1872	1883	1886	rVV2	2151764	4021010	14.01%	0.821%
39	14.093	1886	1888	1901	rVB	1693994	2496974	8.70%	0.510%
40	14.375	1929	1936	1942	rBV	1572680	2555530	8.90%	0.522%
41	14.451	1942	1949	1955	rVB	9569523	16246526	56.59%	3.317%
42	14.504	1955	1958	1965	rVB2	74305	120680	0.42%	0.025%
43	14.634	1966	1980	1986	rBV2	3416167	6149418	21.42%	1.256%
44	14.693	1986	1990	1995	rVB	160402	225411	0.79%	0.046%
45	14.775	1995	2004	2009	rBV	5086224	10332873	35.99%	2.110%
46	15.022	2042	2046	2055	rVB	189823	242459	0.84%	0.050%
47	15.187	2068	2074	2080	rBV2	33240	60066	0.21%	0.012%
48	15.269	2081	2088	2097	rVB3	47658	113355	0.39%	0.023%
49	15.398	2102	2110	2113	rBV	3110653	4703544	16.38%	0.960%
50	15.451	2113	2119	2126	rVB2	15101031	28709848	100.00%	5.862%
51	15.522	2126	2131	2137	rBV	794582	1105077	3.85%	0.226%
52	15.704	2157	2162	2168	rBV	256065	372717	1.30%	0.076%
53	15.957	2195	2205	2215	rBV7	41047	122425	0.43%	0.025%
54	16.140	2230	2236	2245	rBV4	121708	235458	0.82%	0.048%
55	16.363	2262	2274	2282	rBV	3056328	4514092	15.72%	0.922%
56	16.481	2286	2294	2300	rBV	15704347	23545564	82.01%	4.808%
57	16.581	2307	2311	2315	rBV3	37179	53327	0.19%	0.011%
58	16.840	2347	2355	2365	rBV	10762866	17163903	59.78%	3.505%
59	16.945	2370	2373	2376	rVV	62605	97540	0.34%	0.020%
60	16.992	2376	2381	2389	rVB4	98427	211355	0.74%	0.043%
61	17.192	2408	2415	2419	rVV	1650358	2848982	9.92%	0.582%
62	17.240	2419	2423	2428	rVV	4477856	6828924	23.79%	1.394%
63	17.292	2428	2432	2435	rVV	3098948	4746420	16.53%	0.969%
64	17.328	2435	2438	2444	rVB	5115440	6470731	22.54%	1.321%
65	17.498	2464	2467	2476	rVB2	34890	61843	0.22%	0.013%
66	17.616	2482	2487	2492	rVB3	53631	89631	0.31%	0.018%
67	17.816	2517	2521	2526	rVV2	78215	134861	0.47%	0.028%
68	17.892	2530	2534	2541	rVB	298432	391011	1.36%	0.080%
69	18.128	2569	2574	2581	rBV	981995	1468743	5.12%	0.300%
70	18.210	2581	2588	2593	rVV	13008167	19846568	69.13%	4.052%
71	18.504	2634	2638	2648	rVB3	96510	194800	0.68%	0.040%
72	18.598	2650	2654	2658	rVV	48090	75075	0.26%	0.015%
73	18.663	2660	2665	2674	rVB	360152	534419	1.86%	0.109%
74	18.904	2703	2706	2712	rVB	53493	78789	0.27%	0.016%
75	19.039	2725	2729	2734	rBV6	29270	61696	0.21%	0.013%
76	19.116	2739	2742	2745	rVV	107009	123860	0.43%	0.025%

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77	19.216	2755	2759	2762	rBV2	51287	72309	0.25%	0.015%
78	19.257	2762	2766	2769	rVV2	85460	121689	0.42%	0.025%
79	19.328	2770	2778	2788	rVB	12920712	21000484	73.15%	4.288%
80	19.457	2796	2800	2809	rBV	542757	832062	2.90%	0.170%
81	19.575	2816	2820	2824	rBV	125327	172824	0.60%	0.035%
82	19.669	2830	2836	2838	rBV	3461729	4776986	16.64%	0.975%
83	19.698	2838	2841	2853	rVB	6417221	9228162	32.14%	1.884%
84	19.792	2855	2857	2861	rBV3	44096	67020	0.23%	0.014%
85	19.828	2861	2863	2869	rVB3	62389	79152	0.28%	0.016%
86	20.404	2958	2961	2965	rBV	122948	131836	0.46%	0.027%
87	20.657	2999	3004	3009	rBV	3601772	4355217	15.17%	0.889%
88	21.316	3111	3116	3126	rVB	606790	1094212	3.81%	0.223%
89	21.475	3140	3143	3150	rVB2	199488	255700	0.89%	0.052%
90	21.563	3151	3158	3162	rVV	14375997	19541042	68.06%	3.990%
91	21.622	3162	3168	3175	rVV	8647975	18677073	65.05%	3.814%
92	21.704	3175	3182	3191	rVB	12932950	24451611	85.17%	4.993%
93	23.951	3552	3564	3577	rVB	9173286	24028477	83.69%	4.906%
94	24.127	3591	3594	3603	rVB2	173993	359780	1.25%	0.073%
95	24.769	3695	3703	3709	rBV2	1520696	4048308	14.10%	0.827%
96	24.845	3710	3716	3728	rVB	2144016	5522261	19.23%	1.128%
97	24.992	3734	3741	3752	rVB	842253	2302114	8.02%	0.470%
98	28.798	4379	4388	4391	rBV2	538082	1525033	5.31%	0.311%
99	28.904	4394	4406	4407	rBV	2081481	5208606	18.14%	1.064%
100	30.009	4581	4594	4610	rVB	801307	3997737	13.92%	0.816%

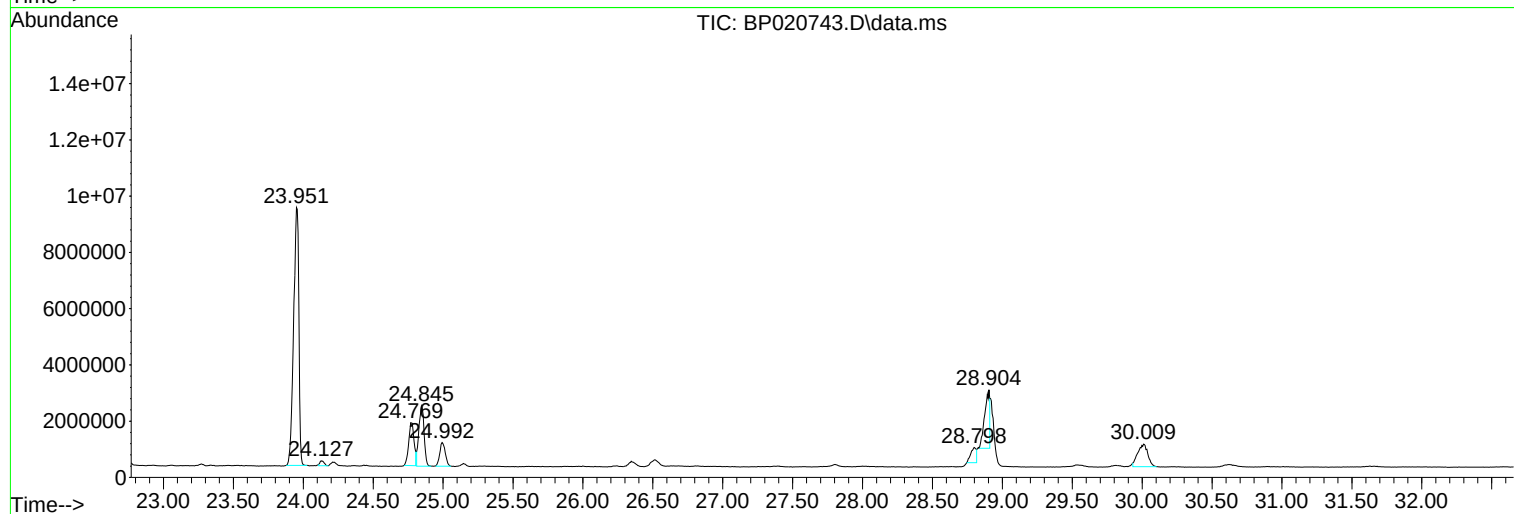
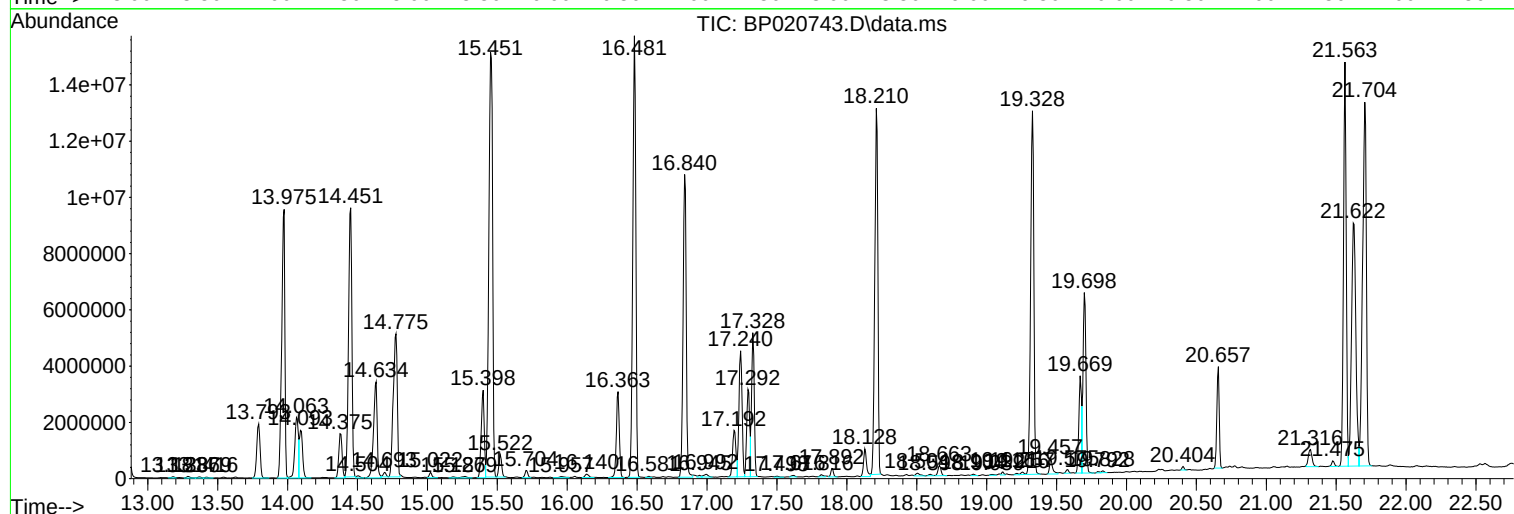
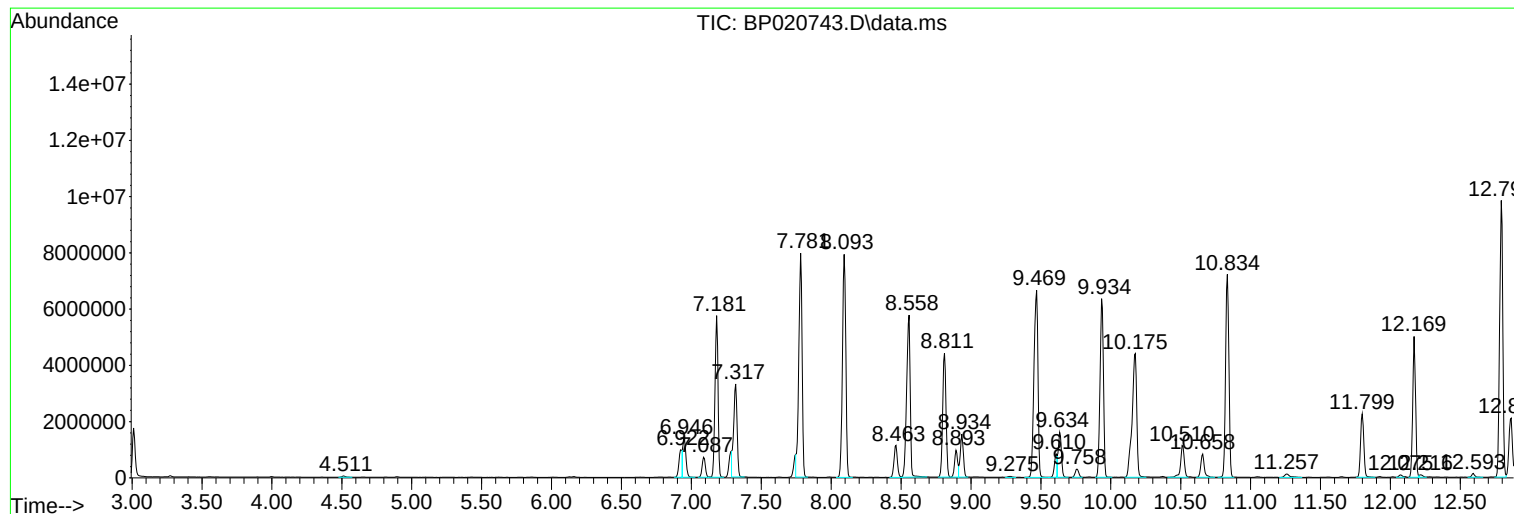
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Instrument :
BNA_P
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062524.MA.M
Quant Title : SVOA CALIBRATION

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



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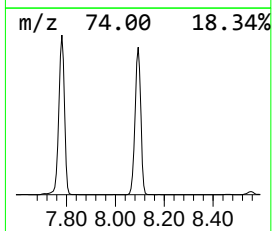
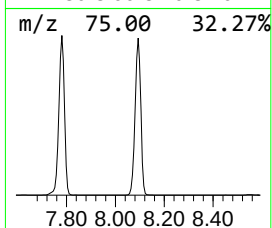
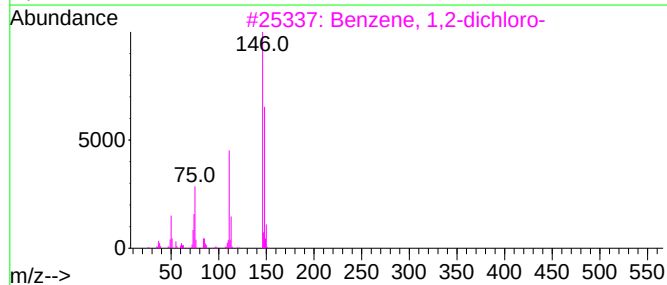
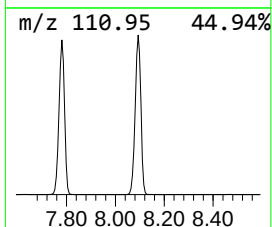
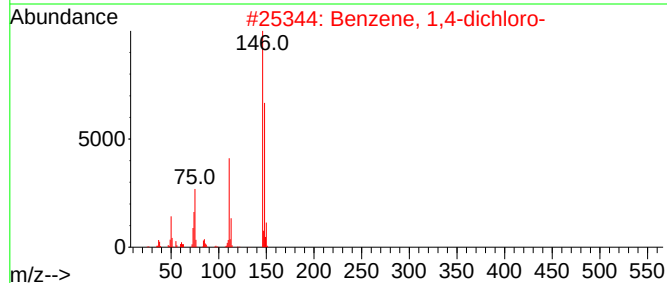
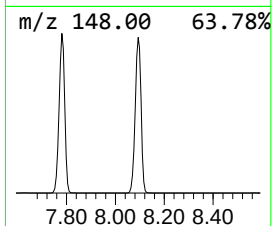
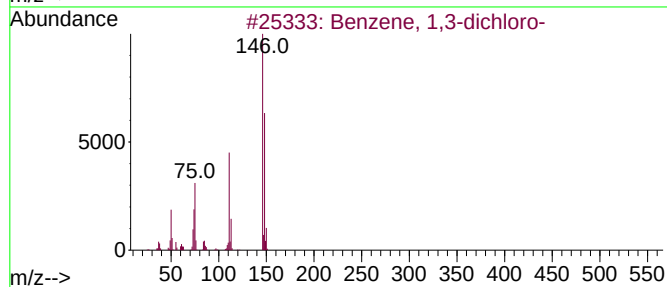
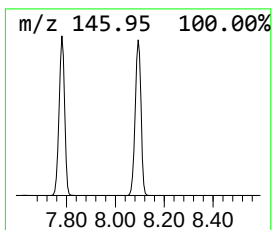
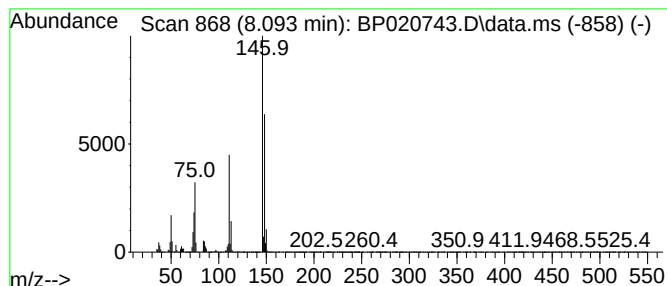
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062524.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Benzene, 1,3-dichloro- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.093	20.04 ng/ul	13090200	1,4-Dichlorobenzene-d4	7.746

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	98
2			Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	98
3			Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	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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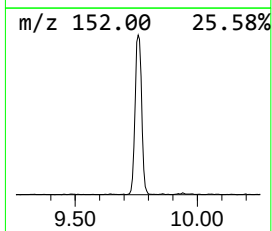
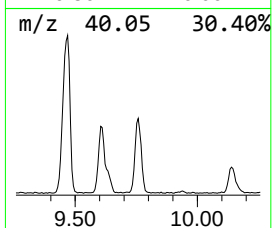
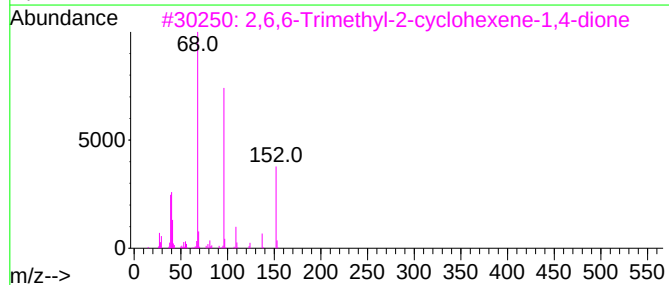
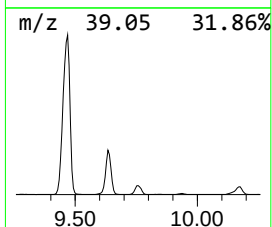
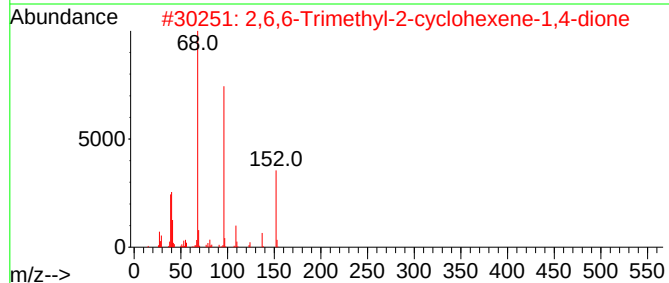
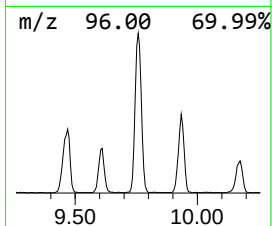
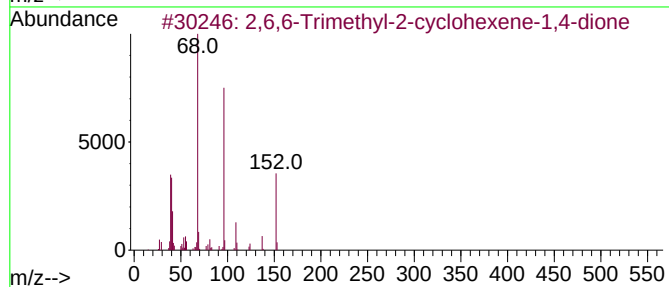
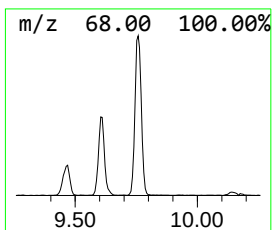
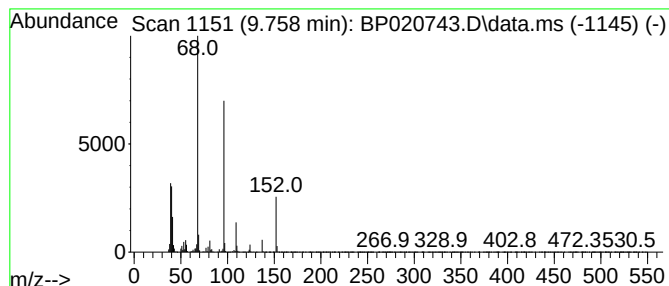
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062524.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 2,6,6-Trimethyl-2-cyclohexe... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.758	5.22 ng/ul	525015	Naphthalene-d8	10.511

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	95		
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	91		
5	2-Cyclohexen-1-one	96	C6H8O	000930-68-7	58		



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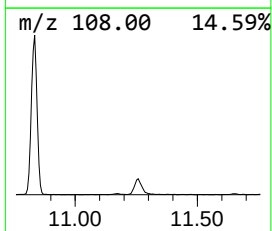
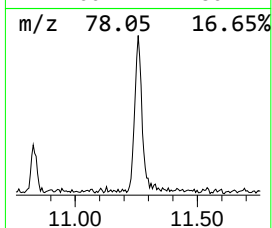
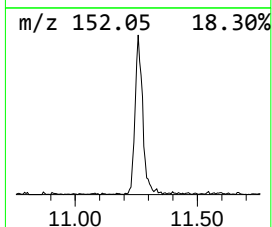
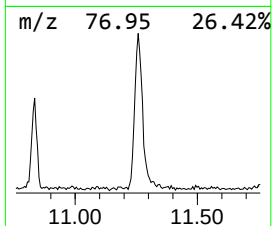
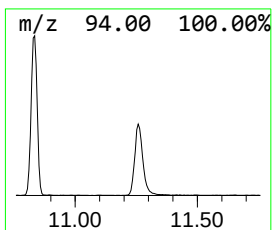
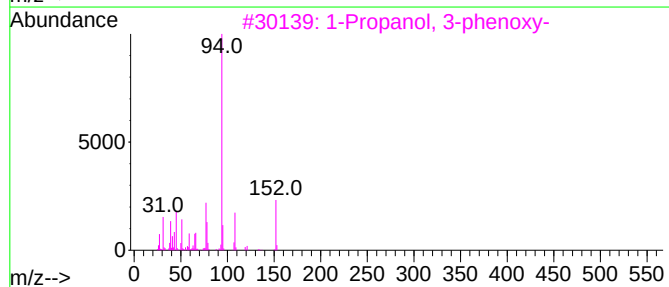
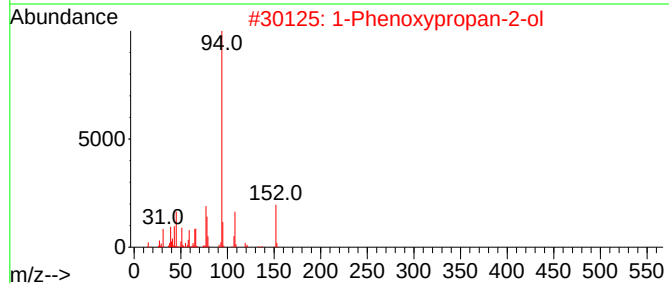
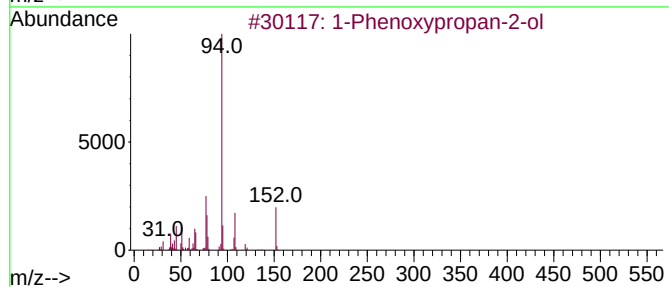
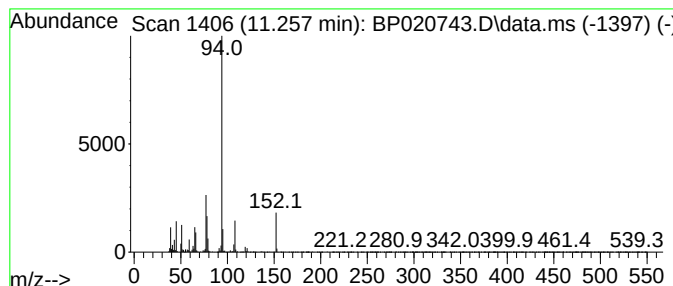
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 1-Phenoxypropan-2-ol Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.258	3.02 ng/ul	303956	Naphthalene-d8	10.511

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	93
2			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	87
3			1-Propanol, 3-phenoxy-	152	C9H12O2	006180-61-6	64
4			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	64
5			1-Propanol, 2-phenoxy-	152	C9H12O2	004169-04-4	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070124\
Data File : BP020743.D
Acq On : 02 Jul 2024 09:26
Operator : MA/JU
Sample : P2963-19
Misc :
ALS Vial : 33 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4DK6

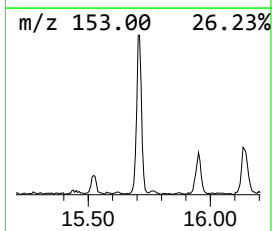
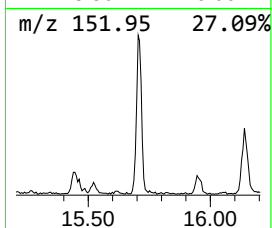
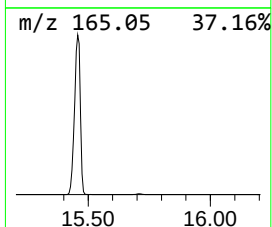
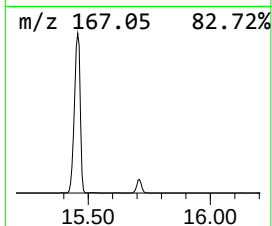
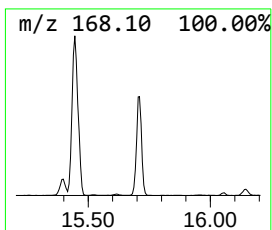
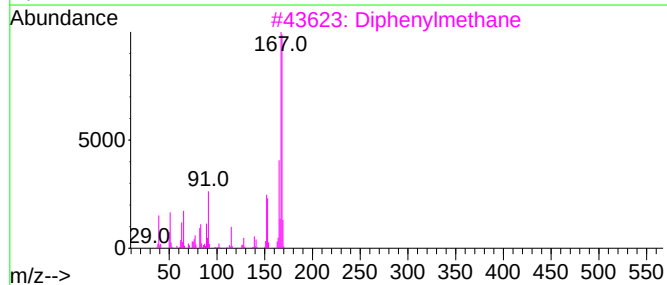
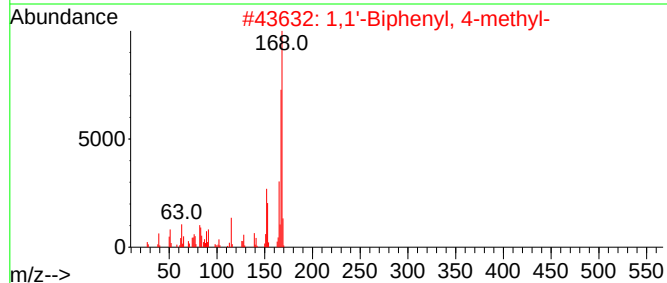
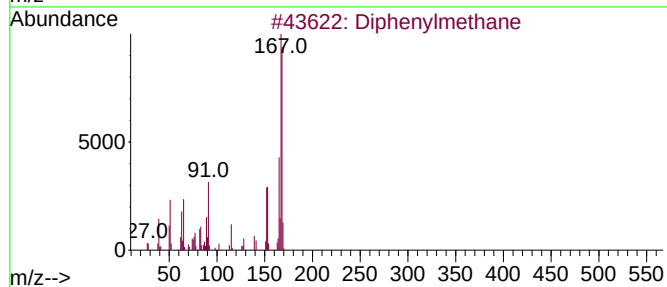
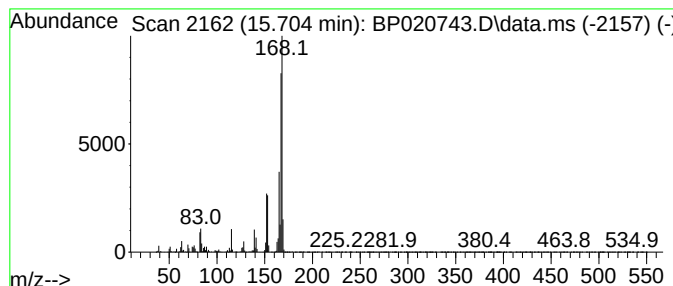
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Diphenylmethane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.704	2.92 ng/ul	372717	Acenaphthene-d10	14.375

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Diphenylmethane	168	C13H12	000101-81-5	93
2			1,1'-Biphenyl, 4-methyl-	168	C13H12	000644-08-6	93
3			Diphenylmethane	168	C13H12	000101-81-5	93
4			1,1'-Biphenyl, 2-methyl-	168	C13H12	000643-58-3	93
5			Diphenylmethane	168	C13H12	000101-81-5	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070124\
Data File : BP020743.D
Acq On : 02 Jul 2024 09:26
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Sample : P2963-19
Misc :
ALS Vial : 33 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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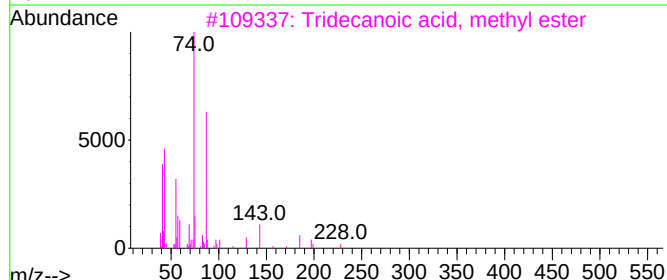
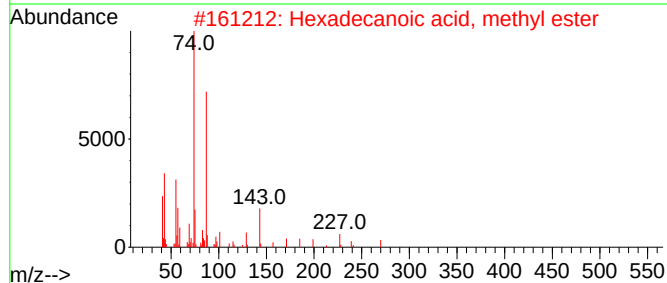
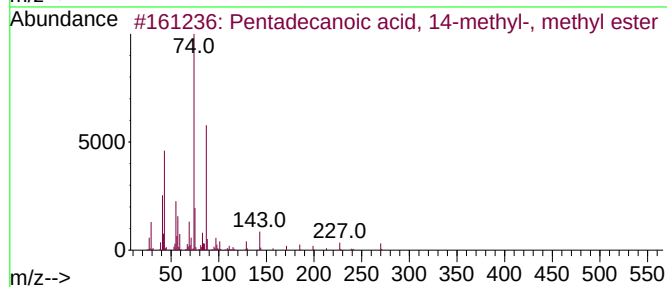
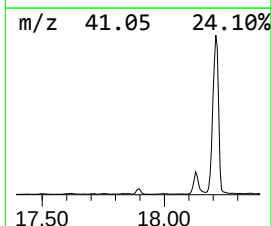
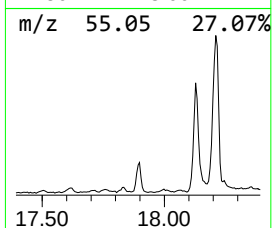
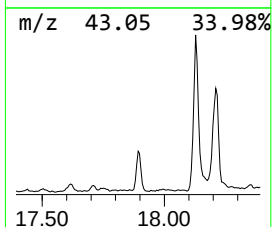
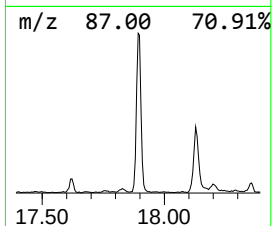
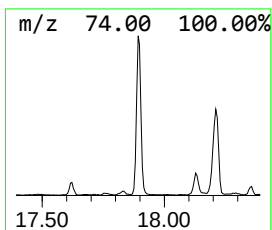
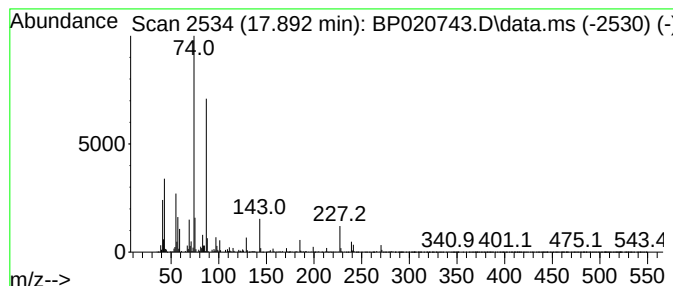
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Pentadecanoic acid, 14-meth... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.892	2.74 ng/ul	391011	Phenanthrene-d10	17.192

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentadecanoic acid, 14-methyl-, ...	270	C17H34O2	005129-60-2	95
2			Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	93
3			Tridecanoic acid, methyl ester	228	C14H28O2	001731-88-0	93
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\BP070124\
Data File : BP020743.D
Acq On : 02 Jul 2024 09:26
Operator : MA/JU
Sample : P2963-19
Misc :
ALS Vial : 33 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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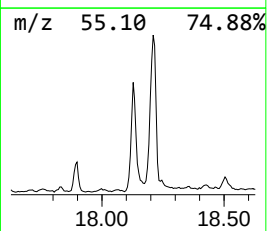
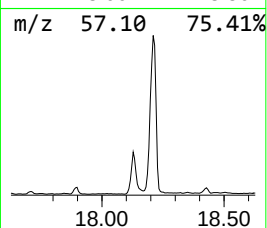
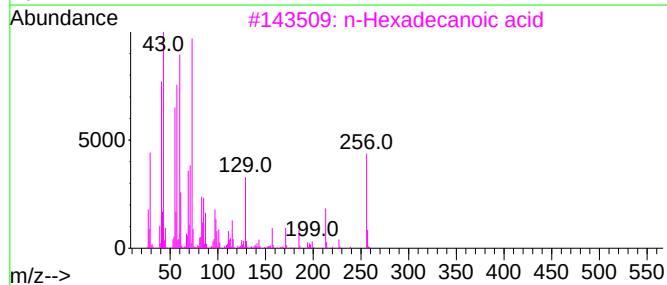
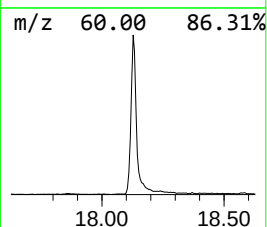
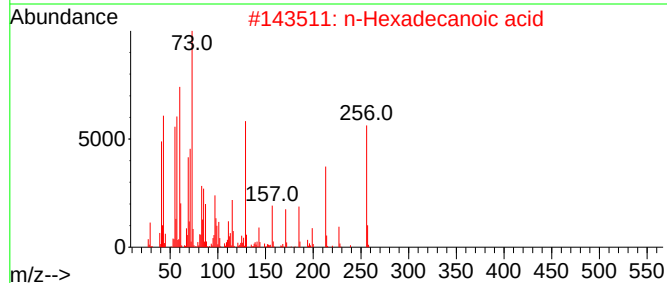
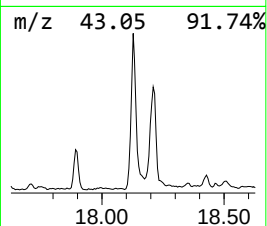
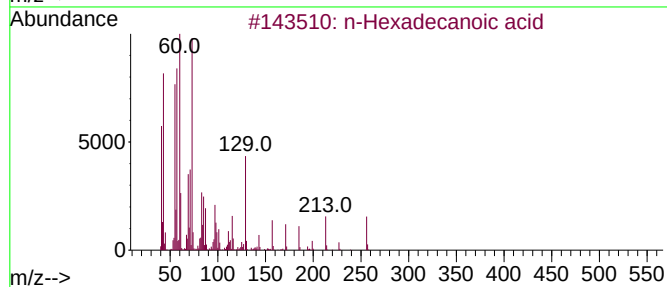
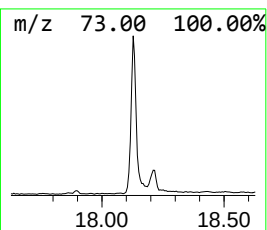
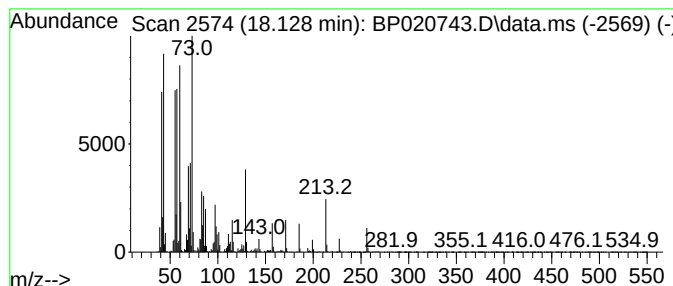
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 n-Hexadecanoic acid Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.128	10.31 ng/ul	1468740	Phenanthrene-d10	17.192

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	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070124\
Data File : BP020743.D
Acq On : 02 Jul 2024 09:26
Operator : MA/JU
Sample : P2963-19
Misc :
ALS Vial : 33 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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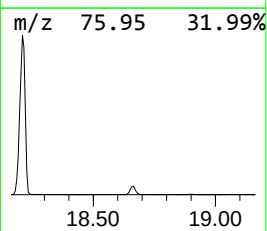
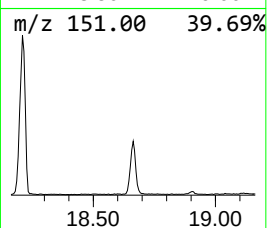
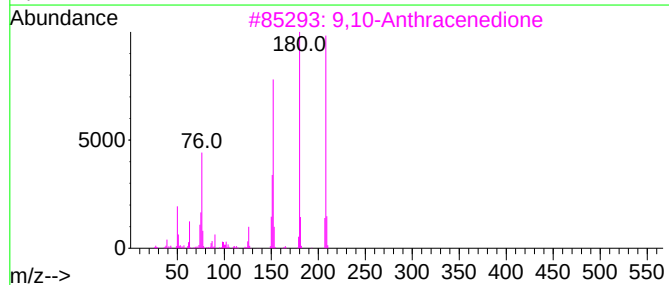
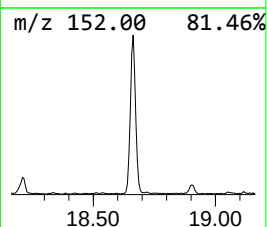
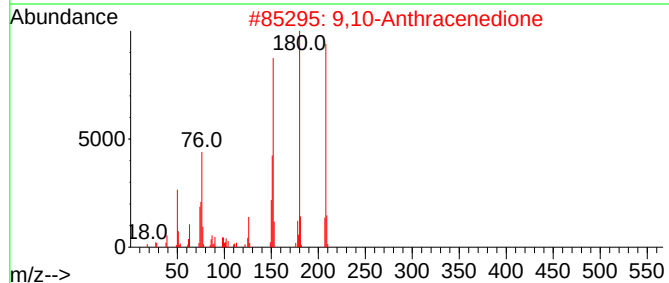
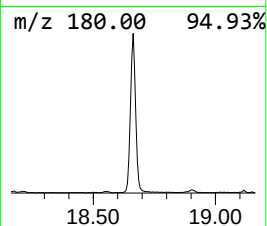
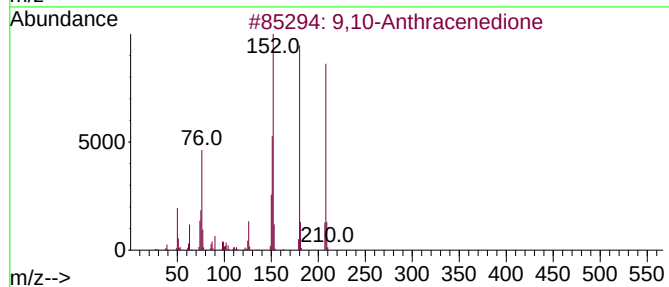
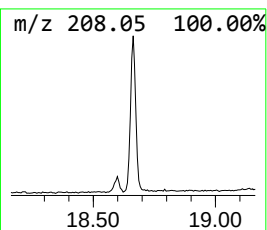
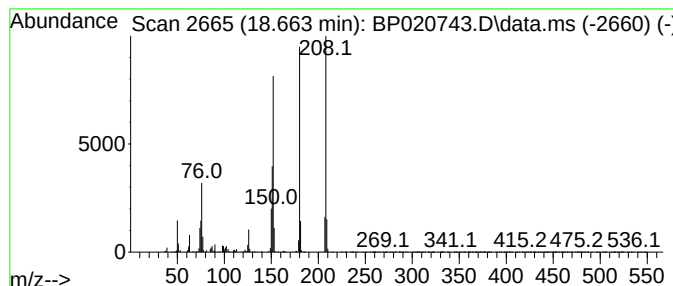
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 9,10-Anthracenedione Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.663	3.75 ng/ul	534419	Phenanthrene-d10	17.192

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	95
4	9,10-Anthracenedione			208	C14H8O2	000084-65-1	93
5	Benzo[c]cinnoline			180	C12H8N2	000230-17-1	70



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070124\
Data File : BP020743.D
Acq On : 02 Jul 2024 09:26
Operator : MA/JU
Sample : P2963-19
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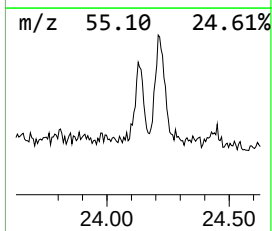
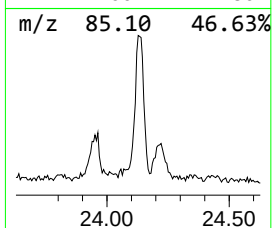
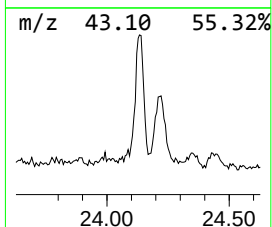
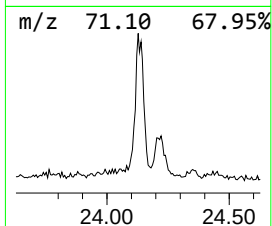
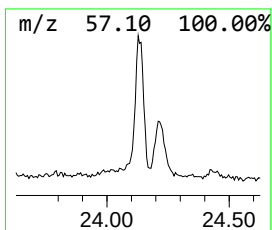
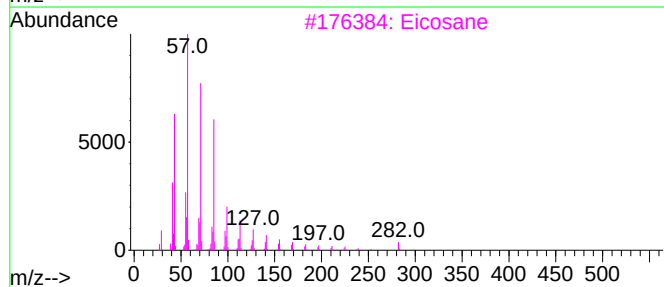
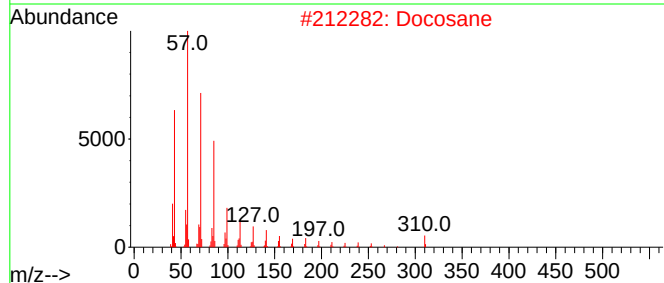
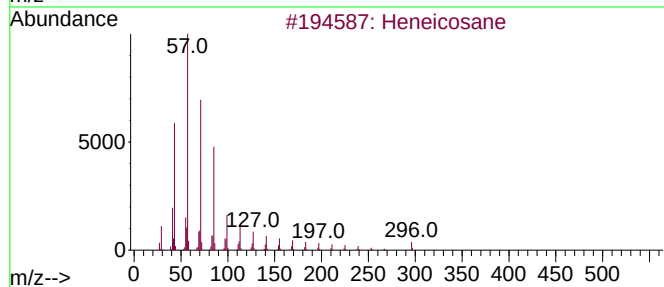
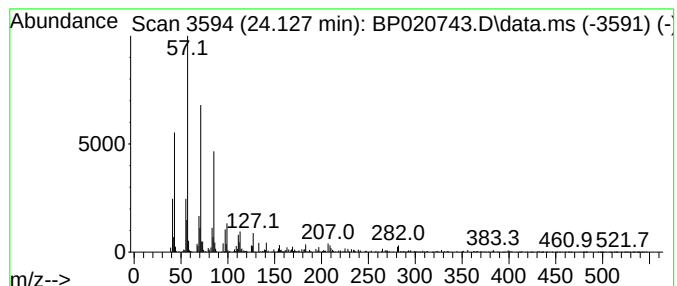
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.127	3.13 ng/ul	359780	Perylene-d12	24.992

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heneicosane	296	C21H44	000629-94-7	97
2			Docosane	310	C22H46	000629-97-0	97
3			Eicosane	282	C20H42	000112-95-8	95
4			Docosane	310	C22H46	000629-97-0	93
5			Tridecane, 1-iodo-	310	C13H27I	035599-77-0	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070124\
Data File : BP020743.D
Acq On : 02 Jul 2024 09:26
Operator : MA/JU
Sample : P2963-19
Misc :
ALS Vial : 33 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062524.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...	9.758	5.2	ng/ul	525015	2	10.511	2012700	20.0
1-Phenoxypropan...	11.258	3.0	ng/ul	303956	2	10.511	2012700	20.0
Diphenylmethane	15.704	2.9	ng/ul	372717	3	14.375	2555530	20.0
Pentadecanoic a...	17.892	2.7	ng/ul	391011	4	17.192	2848980	20.0
n-Hexadecanoic ...	18.128	10.3	ng/ul	1468740	4	17.192	2848980	20.0
9,10-Anthracene...	18.663	3.8	ng/ul	534419	4	17.192	2848980	20.0
(DEL) Alkane: S...	24.127	3.1	ng/ul	359780	6	24.992	2302110	20.0