

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN061424\
 Data File : BN031996.D
 Acq On : 14 Jun 2024 11:41
 Operator : MA/JU
 Sample : P2695-17
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A4DG5

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_N\Methods\SFAM-EPA-BN052924.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BN031996.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.176	28	32	41	rVB	61446	84781	0.26%	0.017%
2	3.446	73	78	86	rBV	36109	60671	0.19%	0.012%
3	3.893	149	154	159	rBV	50068	66329	0.20%	0.013%
4	6.034	512	518	521	rBV	51073	79120	0.24%	0.016%
5	6.864	646	659	672	rVB2	2550521	5805158	17.77%	1.144%
6	6.999	675	682	691	rBV	971347	1537138	4.71%	0.303%
7	7.093	691	698	707	rVB	3908986	5882330	18.01%	1.159%
8	7.193	707	715	717	rBV	1202432	2008521	6.15%	0.396%
9	7.222	717	720	727	rVB	2131973	3212686	9.83%	0.633%
10	7.540	767	774	784	rBV	3435939	5595589	17.13%	1.103%
11	7.658	784	794	796	rVV	1043267	1800634	5.51%	0.355%
12	7.693	796	800	812	rVV	3003934	4793451	14.67%	0.945%
13	7.905	831	836	844	rBV2	34700	63141	0.19%	0.012%
14	8.005	845	853	871	rVB	9928431	17075052	52.27%	3.365%
15	8.381	909	917	924	rBV	1498316	2455794	7.52%	0.484%
16	8.475	924	933	942	rBV	5935142	10463637	32.03%	2.062%
17	8.816	983	991	994	rBV	1099676	1944085	5.95%	0.383%
18	8.863	994	999	1006	rVB	4688779	7573579	23.18%	1.493%
19	9.216	1053	1059	1065	rVB2	50327	89264	0.27%	0.018%
20	9.563	1104	1118	1130	rBV2	4325612	8986131	27.51%	1.771%
21	9.863	1161	1169	1180	rVB	4380603	7293696	22.33%	1.437%
22	10.093	1196	1208	1222	rBV2	7095158	14361099	43.96%	2.830%
23	10.299	1236	1243	1251	rVB3	32255	57940	0.18%	0.011%
24	10.440	1259	1267	1275	rVB	1393195	2348422	7.19%	0.463%
25	10.581	1285	1291	1311	rVB	681269	1241698	3.80%	0.245%
26	10.734	1311	1317	1326	rBV	39266	67205	0.21%	0.013%
27	11.010	1358	1364	1369	rBV2	84564	139727	0.43%	0.028%
28	11.063	1369	1373	1380	rVB2	81109	123882	0.38%	0.024%
29	11.187	1388	1394	1401	rBV	199321	388867	1.19%	0.077%
30	11.740	1478	1488	1501	rBV	7359007	12437865	38.08%	2.451%
31	12.534	1618	1623	1629	rBV	83517	117218	0.36%	0.023%
32	12.722	1648	1655	1661	rVV	3949899	6317513	19.34%	1.245%
33	12.799	1661	1668	1683	rVV	4999643	8341406	25.54%	1.644%
34	12.928	1685	1690	1693	rVV6	32430	72952	0.22%	0.014%
35	12.963	1693	1696	1701	rVB	49026	69794	0.21%	0.014%
36	13.716	1817	1824	1833	rBV	2220565	3205017	9.81%	0.632%

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Smoothing : OFF

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Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

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Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN052924.MA.M
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37	13.893	1846	1854	1861	rBV	4820142	7006390	21.45%	1.381%
38	13.993	1861	1871	1881	rVB	2755733	4360085	13.35%	0.859%
39	14.304	1916	1924	1929	rBV	1868535	2982483	9.13%	0.588%
40	14.369	1929	1935	1941	rVB	13059858	21399475	65.51%	4.217%

41	14.428	1941	1945	1951	rVB	258118	348014	1.07%	0.069%
42	14.546	1953	1965	1975	rBV3	6838147	16211440	49.63%	3.195%
43	14.693	1982	1990	1995	rBV	7285009	11417579	34.95%	2.250%
44	14.728	1995	1996	2006	rVV2	116601	133870	0.41%	0.026%
45	14.940	2027	2032	2040	rVB	117034	187487	0.57%	0.037%

46	15.134	2059	2065	2071	rBV	5167007	7094637	21.72%	1.398%
47	15.175	2071	2072	2085	rVB4	62724	90698	0.28%	0.018%
48	15.298	2086	2093	2098	rBV	3287921	5016528	15.36%	0.989%
49	15.357	2098	2103	2110	rVB	10013539	15197949	46.52%	2.995%
50	15.434	2110	2116	2127	rBV	895955	1294666	3.96%	0.255%

51	15.604	2140	2145	2150	rBV	191682	261893	0.80%	0.052%
52	15.851	2181	2187	2194	rBV2	78752	160346	0.49%	0.032%
53	16.028	2211	2217	2224	rBV3	135037	250342	0.77%	0.049%
54	16.245	2242	2254	2261	rVB	5496008	7726173	23.65%	1.523%
55	16.463	2286	2291	2294	rBV2	55451	87147	0.27%	0.017%

56	16.704	2325	2332	2342	rBV	8159120	12559789	38.45%	2.475%
57	16.857	2353	2358	2369	rVB3	151110	304210	0.93%	0.060%
58	16.992	2372	2381	2384	rBV4	35843	74182	0.23%	0.015%
59	17.051	2385	2391	2394	rVV	2053652	3140648	9.61%	0.619%
60	17.098	2394	2399	2404	rVV	10886901	16568760	50.72%	3.265%

61	17.151	2404	2408	2411	rVV	3600145	5247486	16.06%	1.034%
62	17.187	2411	2414	2420	rVB	4553529	5823001	17.83%	1.148%
63	17.340	2436	2440	2447	rVB2	42890	66783	0.20%	0.013%
64	17.463	2453	2461	2466	rVB4	89092	168551	0.52%	0.033%
65	17.598	2480	2484	2487	rVV4	40508	68814	0.21%	0.014%

66	17.663	2491	2495	2501	rVV2	104331	179680	0.55%	0.035%
67	17.728	2501	2506	2512	rVB	479600	618608	1.89%	0.122%
68	17.957	2539	2545	2551	rBV	2673612	3694823	11.31%	0.728%
69	18.028	2551	2557	2562	rVB	13515855	19335484	59.19%	3.811%
70	18.239	2589	2593	2597	rBV3	51428	90462	0.28%	0.018%

71	18.322	2603	2607	2616	rVB2	171668	296268	0.91%	0.058%
72	18.416	2619	2623	2629	rVV	60837	105741	0.32%	0.021%
73	18.481	2629	2634	2648	rVB2	492419	802453	2.46%	0.158%
74	18.822	2689	2692	2697	rBV3	64867	115948	0.35%	0.023%
75	18.875	2698	2701	2703	rVV2	69275	80909	0.25%	0.016%

76	18.904	2703	2706	2710	rVB	361358	430269	1.32%	0.085%
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77	19.010	2721	2724	2726	rBV2	66974	81955	0.25%	0.016%
78	19.045	2726	2730	2735	rVV2	214210	374286	1.15%	0.074%
79	19.122	2735	2743	2755	rVB	9474405	13930870	42.65%	2.746%
80	19.239	2759	2763	2778	rBV	1736734	2876642	8.81%	0.567%
81	19.363	2779	2784	2789	rVB	474673	681892	2.09%	0.134%
82	19.492	2794	2806	2811	rVV2	16992014	32666333	100.00%	6.438%
83	19.563	2815	2818	2823	rBV	123075	173559	0.53%	0.034%
84	19.981	2885	2889	2890	rBV	131385	167616	0.51%	0.033%
85	20.145	2914	2917	2921	rBV	247068	256182	0.78%	0.050%
86	20.333	2947	2949	2952	rVB	106978	101681	0.31%	0.020%
87	20.392	2954	2959	2964	rVB	10817781	12731963	38.98%	2.509%
88	20.975	3054	3058	3066	rBV2	914474	1341024	4.11%	0.264%
89	21.122	3080	3083	3088	rVB	467590	540205	1.65%	0.106%
90	21.198	3091	3096	3102	rVV	13462504	18835835	57.66%	3.712%
91	21.280	3103	3110	3115	rVV2	9080492	16536862	50.62%	3.259%
92	21.339	3115	3120	3127	rVB	5939436	8958529	27.42%	1.766%
93	22.222	3266	3270	3276	rVB	380304	588525	1.80%	0.116%
94	22.310	3278	3285	3293	rVB	13127930	23071559	70.63%	4.547%
95	23.298	3445	3453	3458	rBV	3529518	8243298	25.23%	1.625%
96	23.363	3458	3464	3473	rVB	6095059	13407972	41.05%	2.642%
97	24.022	3568	3576	3582	rBV2	2456657	6088932	18.64%	1.200%
98	24.086	3582	3587	3599	rVB	2621421	6026214	18.45%	1.188%
99	24.222	3603	3610	3620	rVB	1602689	3766965	11.53%	0.742%
100	27.604	4161	4185	4202	rVB2	6122287	28788432	88.13%	5.674%

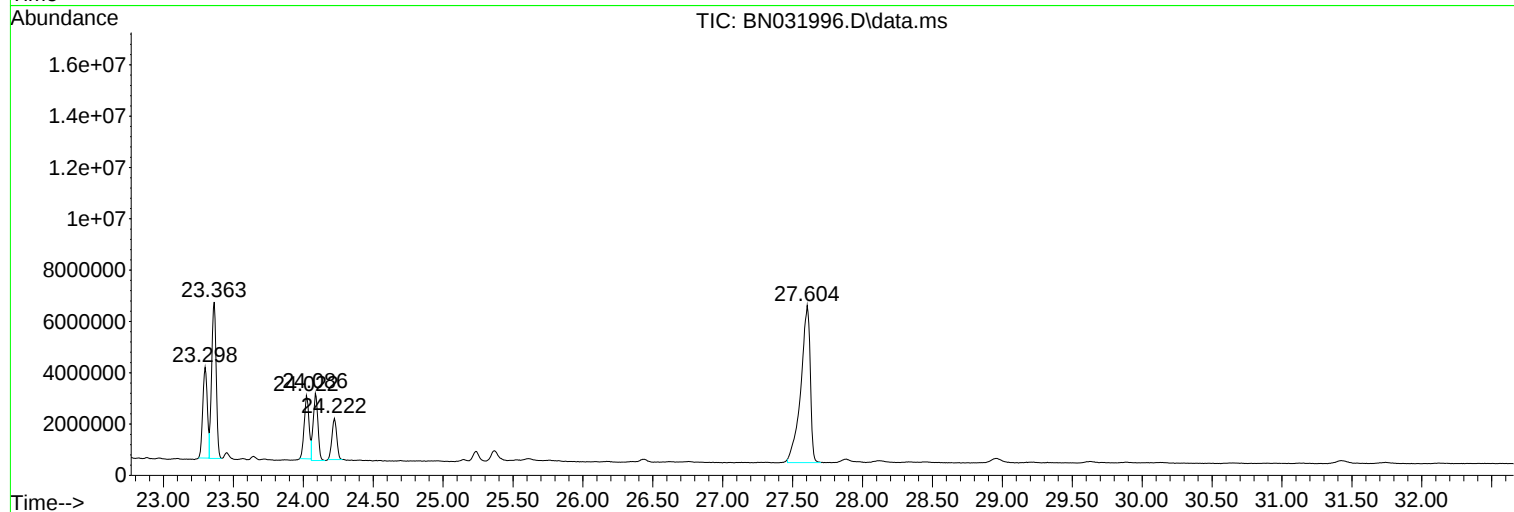
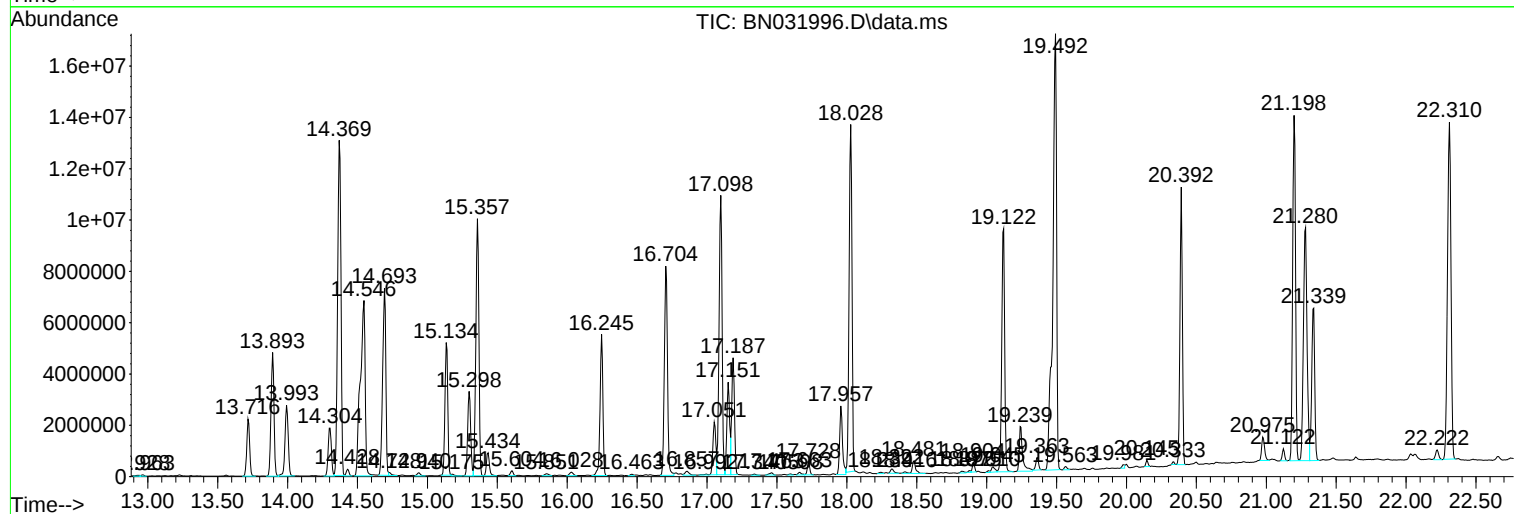
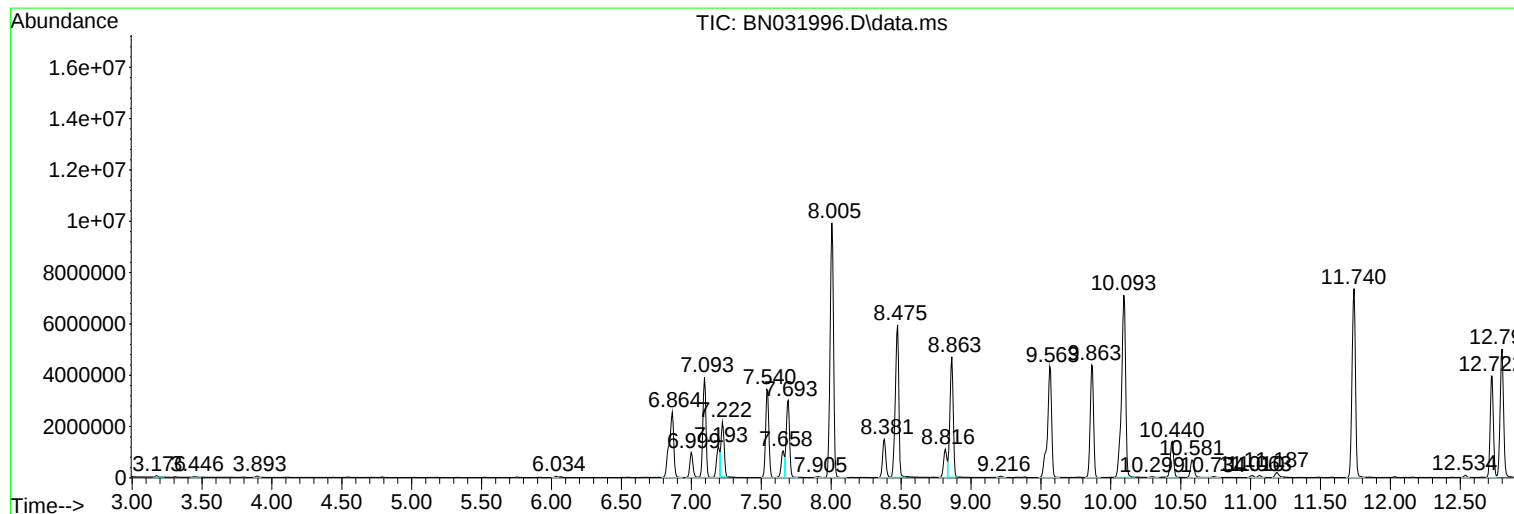
Sum of corrected areas: 507398794

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Instrument :
BNA_N
ClientSampleId :
A4DG5

Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN052924.MA.M
Quant Title : SVOA CALIBRATION

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



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Instrument :
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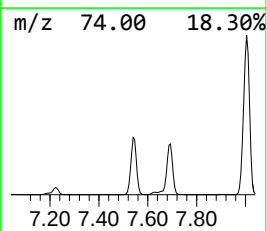
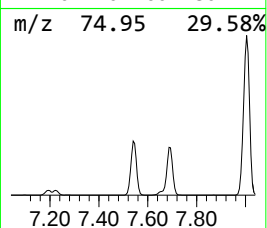
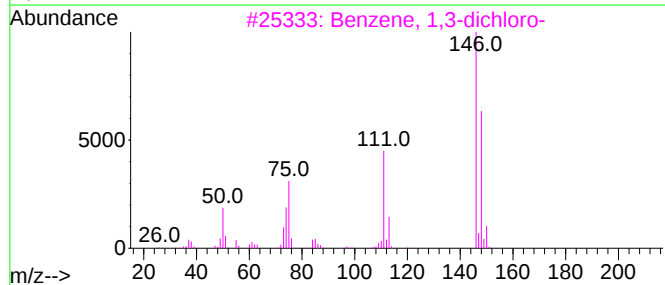
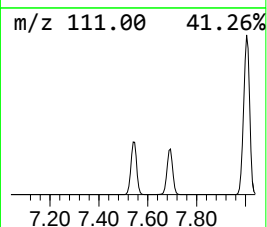
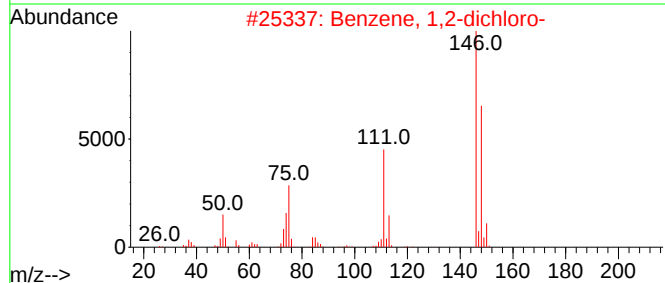
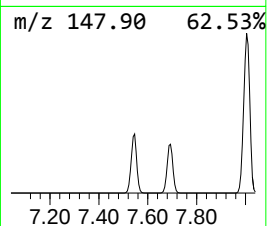
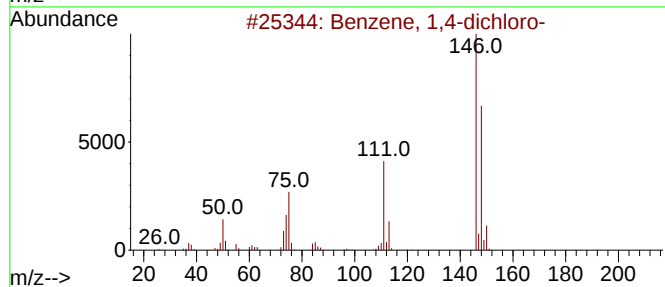
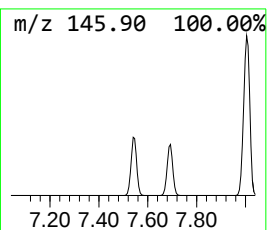
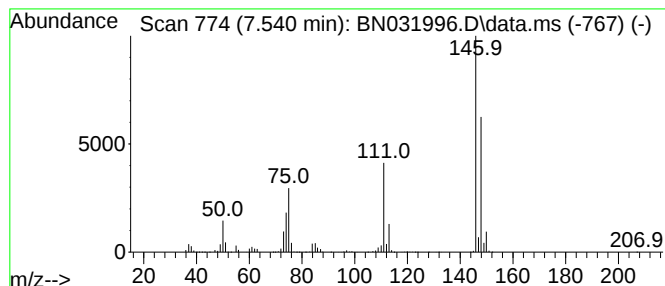
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN052924.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Benzene, 1,4-dichloro- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.540	62.15 ng/ul	5595590	1,4-Dichlorobenzene-d4	7.658

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



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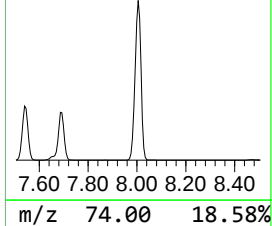
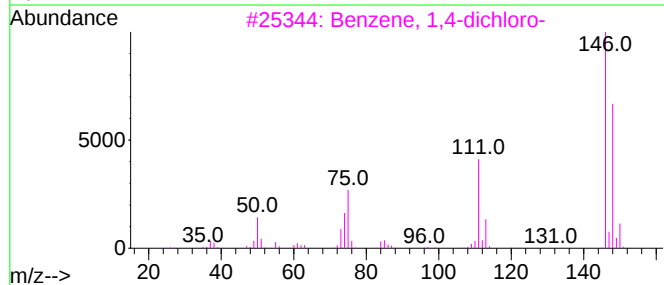
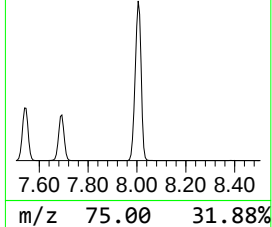
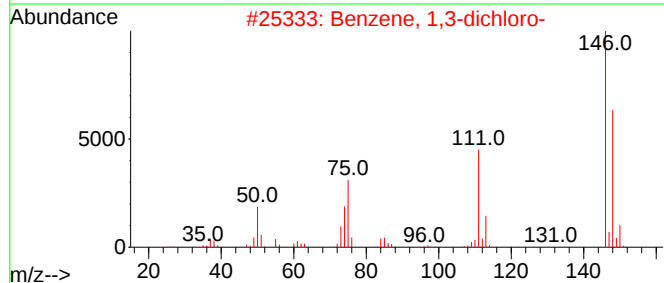
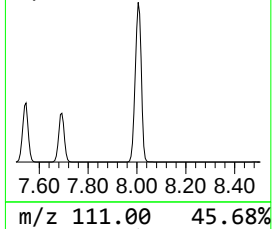
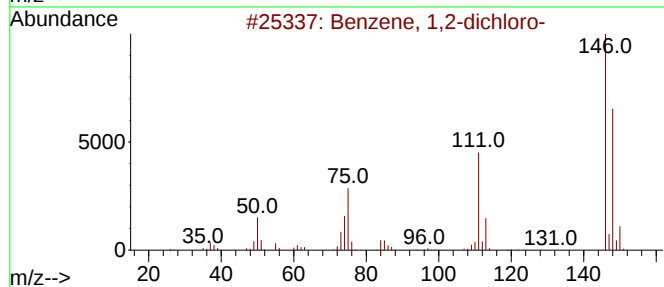
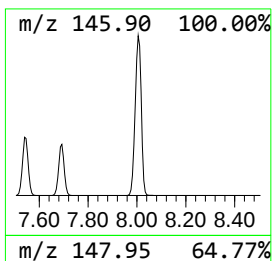
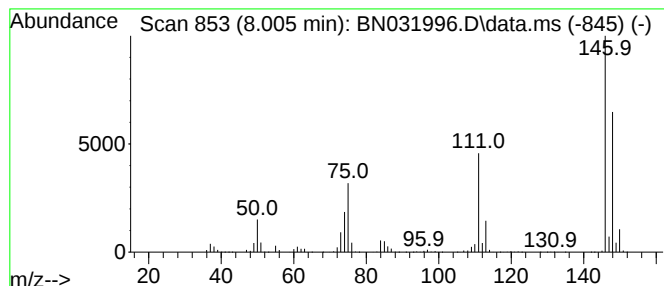
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN052924.MA.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.005	189.66 ng/ul	17075100	1,4-Dichlorobenzene-d4	7.658

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	98
2			Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	98
3			Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	98
4			Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	97
5			Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	96



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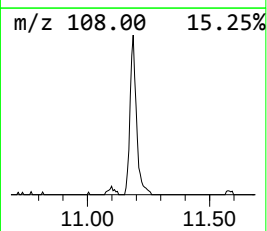
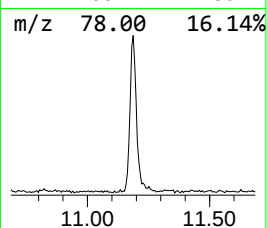
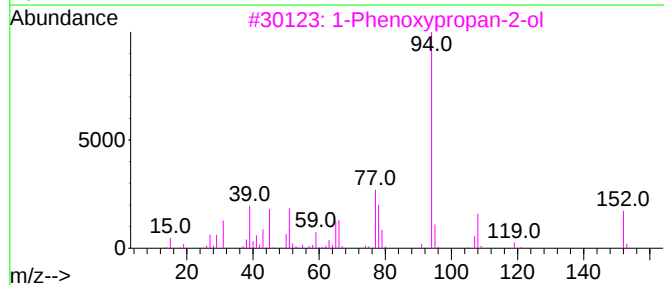
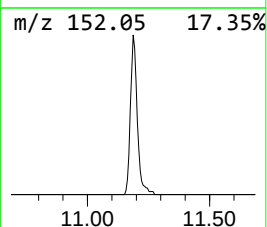
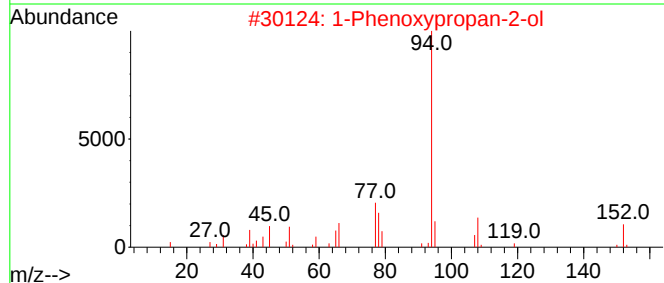
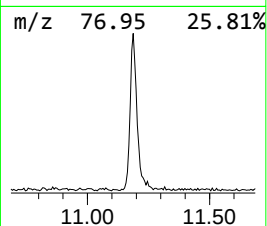
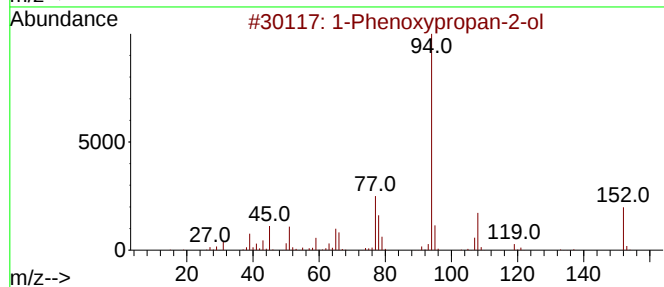
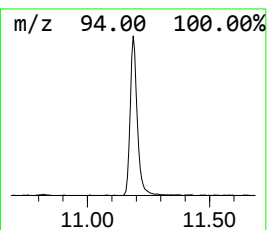
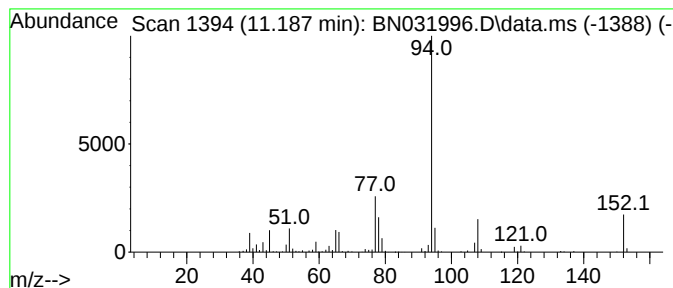
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 1-Phenoxypropan-2-ol Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.187	3.31 ng/ul	388867	Naphthalene-d8	10.440

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	94
3			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	94
4			1-Propanol, 3-phenoxy-	152	C9H12O2	006180-61-6	83
5			Ethanol, 2-phenoxy-	138	C8H10O2	000122-99-6	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN061424\
Data File : BN031996.D
Acq On : 14 Jun 2024 11:41
Operator : MA/JU
Sample : P2695-17
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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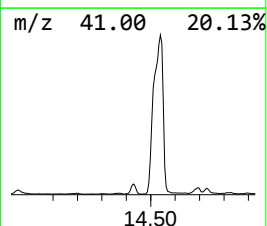
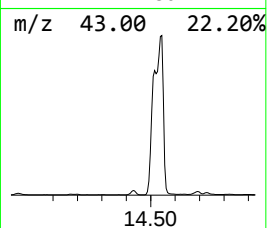
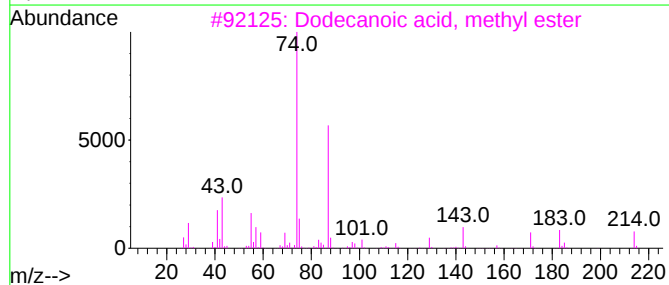
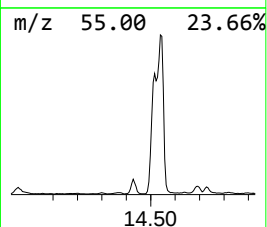
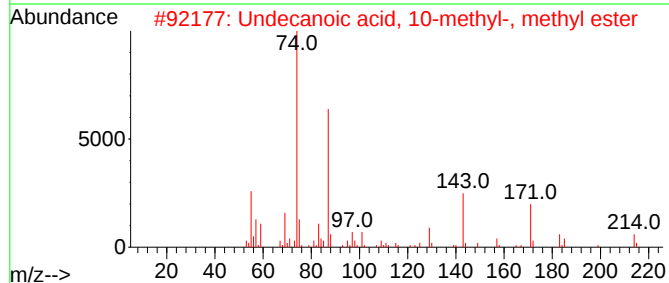
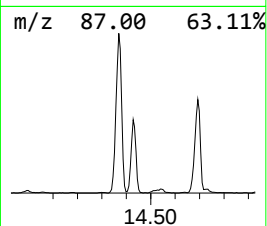
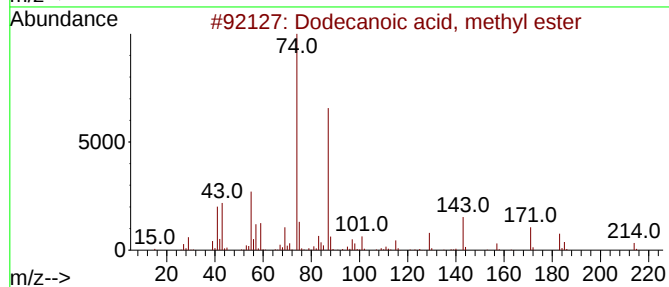
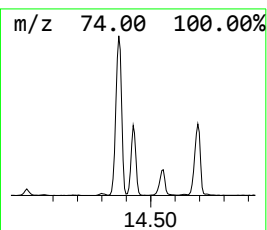
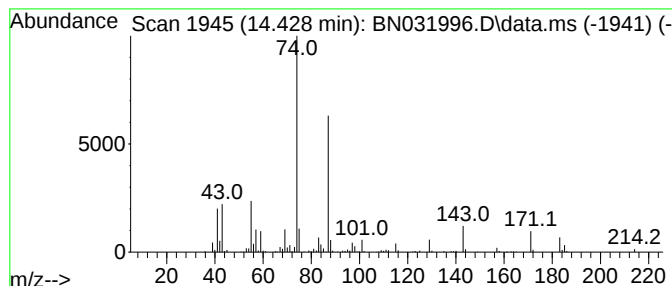
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Dodecanoic acid, methyl ester Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.428	2.33 ng/ul	348014	Acenaphthene-d10	14.304

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dodecanoic acid, methyl ester	214	C13H26O2	000111-82-0	97
2			Undecanoic acid, 10-methyl-, met...	214	C13H26O2	005129-56-6	92
3			Dodecanoic acid, methyl ester	214	C13H26O2	000111-82-0	91
4			Dodecanoic acid, methyl ester	214	C13H26O2	000111-82-0	91
5			Dodecanoic acid, methyl ester	214	C13H26O2	000111-82-0	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN061424\
Data File : BN031996.D
Acq On : 14 Jun 2024 11:41
Operator : MA/JU
Sample : P2695-17
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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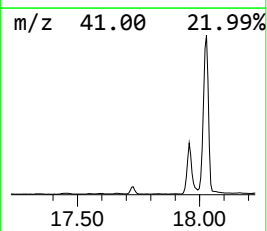
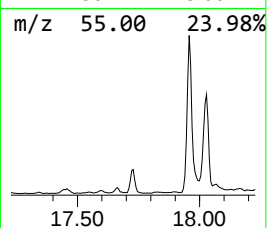
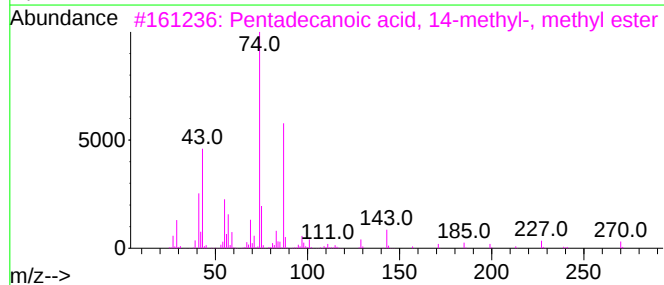
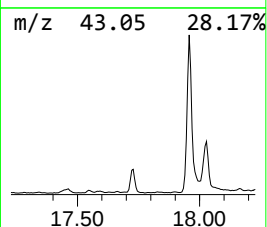
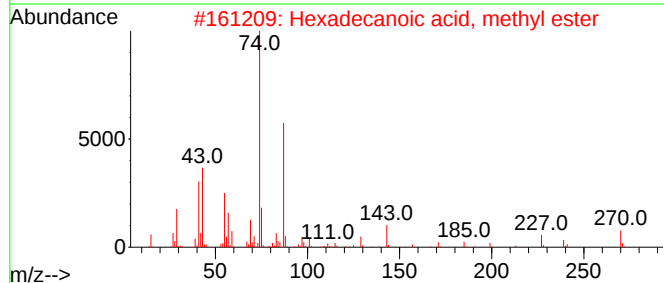
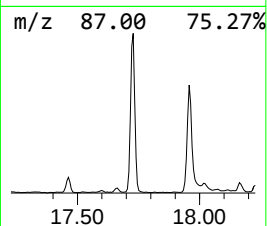
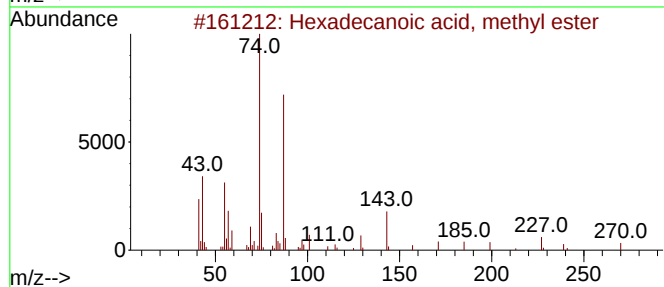
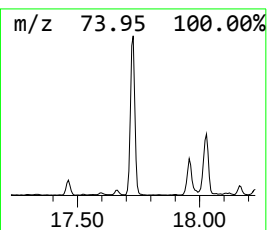
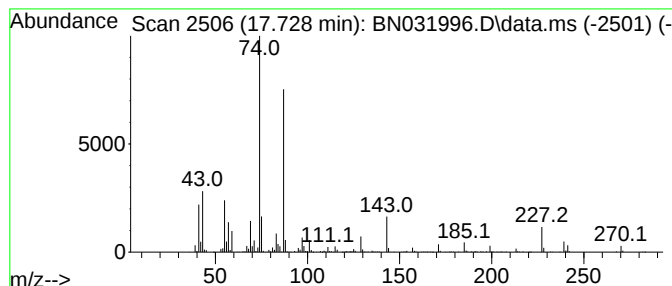
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.728	3.94 ng/ul	618608	Phenanthrene-d10	17.051

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	98
2			Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	97
3			Pentadecanoic acid, 14-methyl-, ...	270	C17H34O2	005129-60-2	96
4			Pentadecanoic acid, 14-methyl-, ...	270	C17H34O2	005129-60-2	90
5			Tridecanoic acid, methyl ester	228	C14H28O2	001731-88-0	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN061424\
Data File : BN031996.D
Acq On : 14 Jun 2024 11:41
Operator : MA/JU
Sample : P2695-17
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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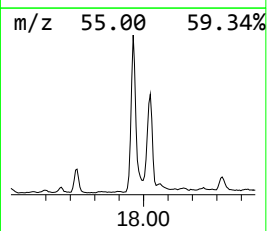
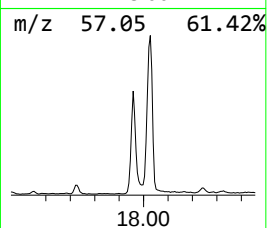
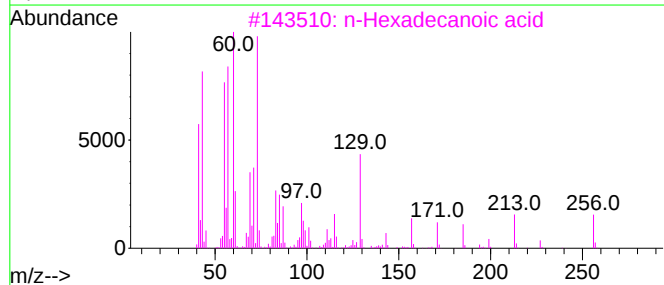
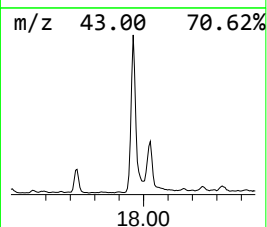
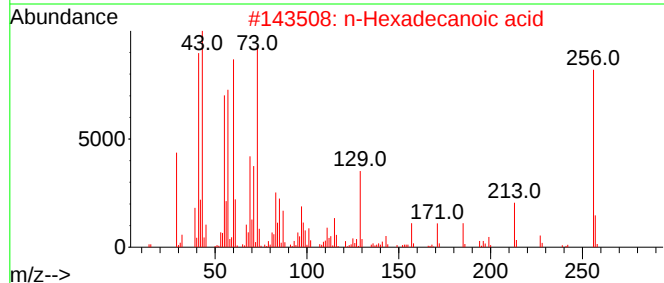
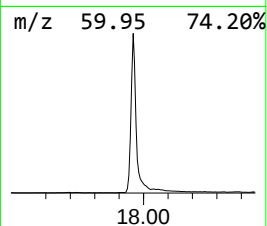
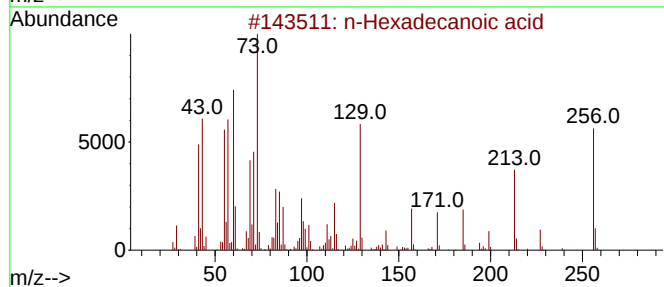
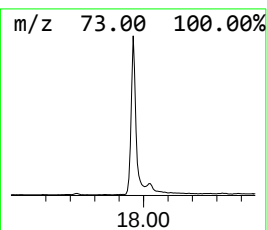
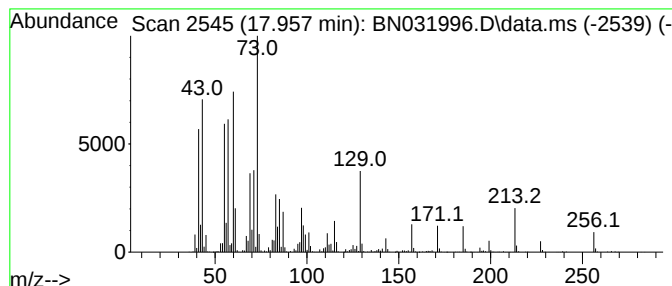
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.957	23.53 ng/ul	3694820	Phenanthrene-d10	17.051

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	99		
4	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98		
5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	86		



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN061424\
Data File : BN031996.D
Acq On : 14 Jun 2024 11:41
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Sample : P2695-17
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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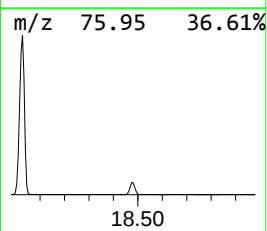
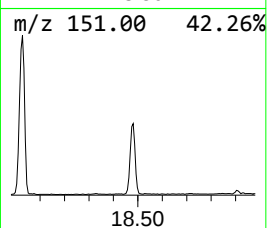
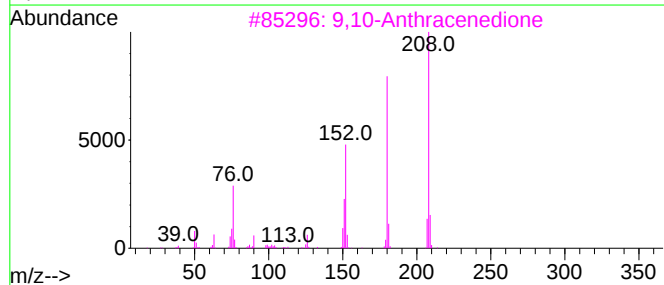
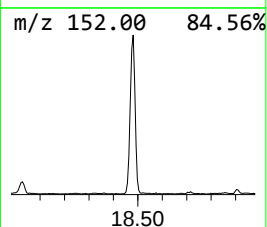
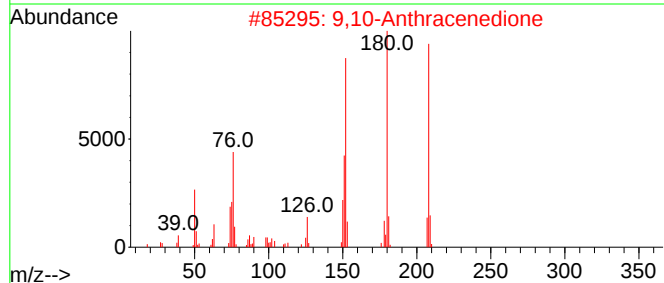
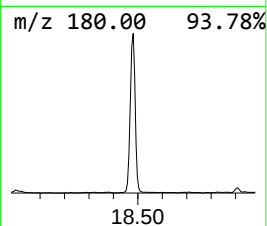
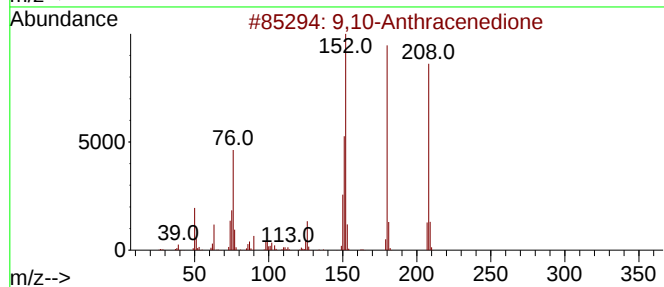
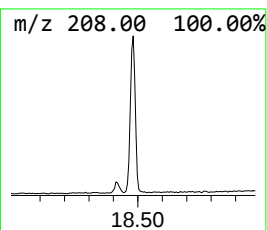
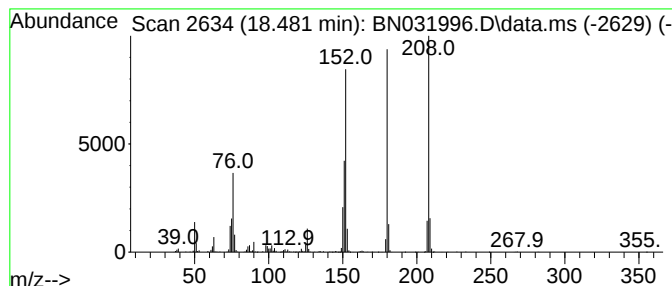
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Quant Title : SVOA CALIBRATION

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.
18.481	5.11 ng/ul	802453	Phenanthrene-d10	17.051

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	94
5	Benzo[c]cinnoline			180	C12H8N2	000230-17-1	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN061424\
Data File : BN031996.D
Acq On : 14 Jun 2024 11:41
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Sample : P2695-17
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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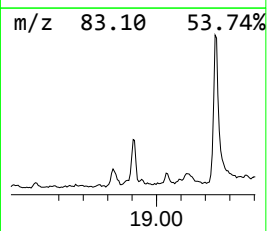
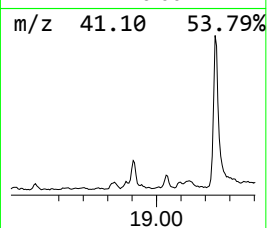
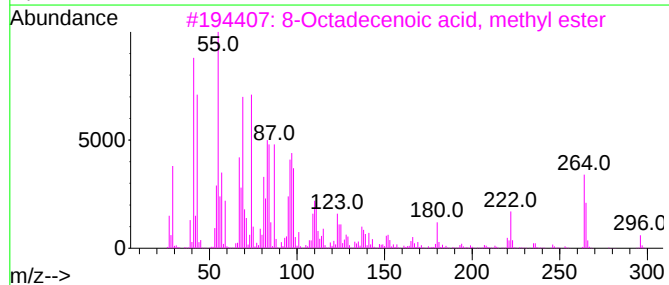
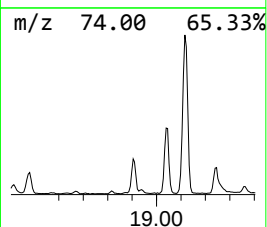
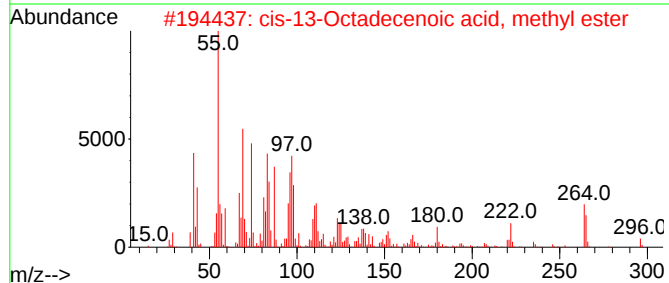
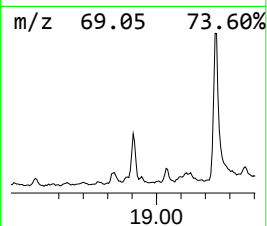
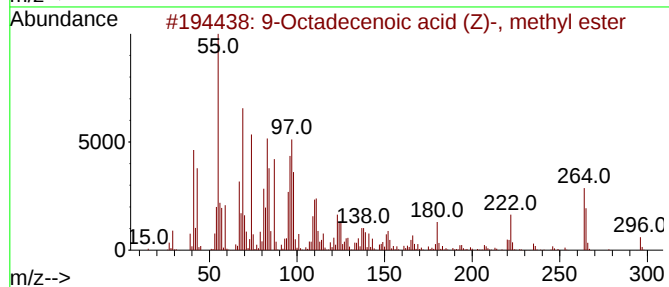
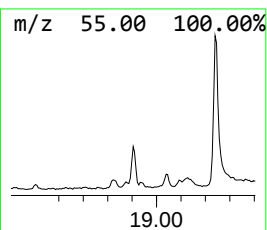
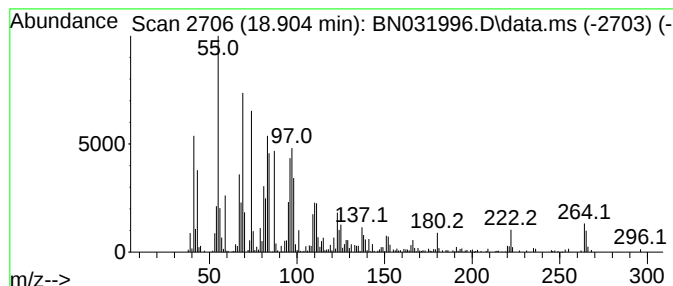
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.904	2.74 ng/ul	430269	Phenanthrene-d10	17.051

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			9-Octadecenoic acid (Z)-, methyl...	296	C19H36O2	000112-62-9	99
2			cis-13-Octadecenoic acid, methyl...	296	C19H36O2	1010333-58-3	99
3			8-Octadecenoic acid, methyl ester	296	C19H36O2	002345-29-1	99
4			trans-13-Octadecenoic acid, meth...	296	C19H36O2	1000333-61-3	99
5			6-Octadecenoic acid, methyl ester	296	C19H36O2	052355-31-4	98



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN061424\
Data File : BN031996.D
Acq On : 14 Jun 2024 11:41
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Sample : P2695-17
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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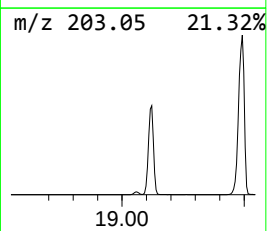
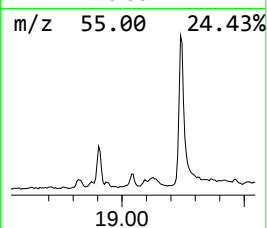
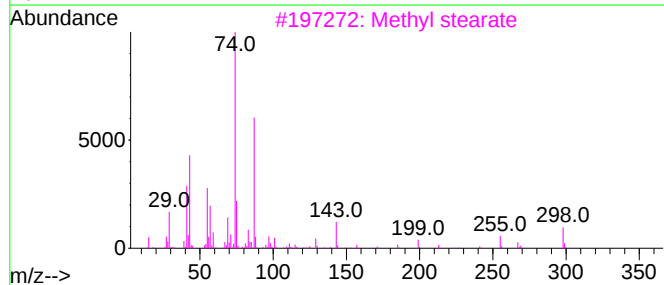
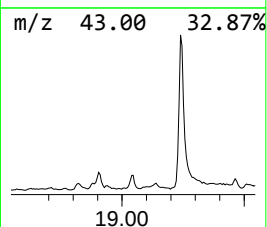
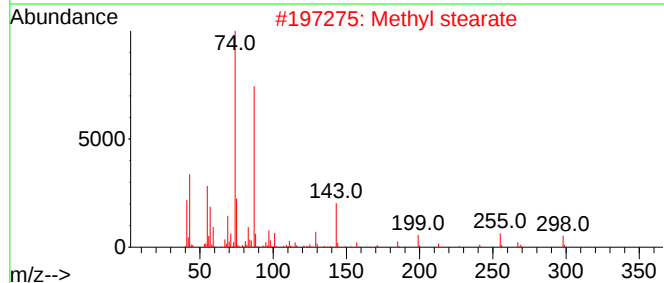
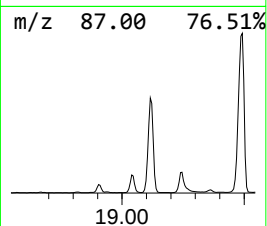
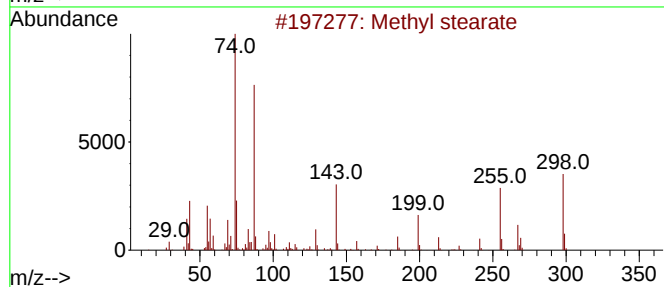
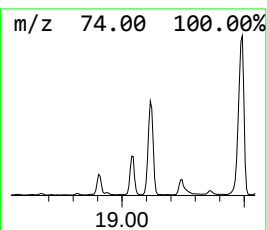
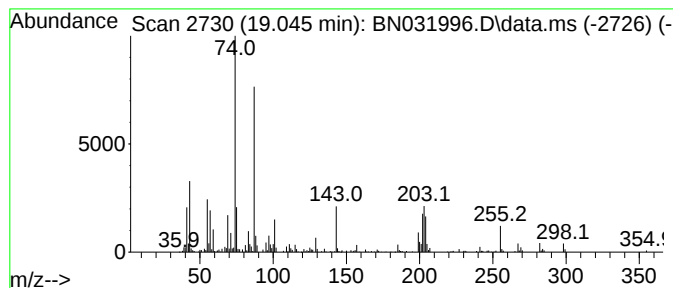
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 Methyl stearate Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.045	2.38 ng/ul	374286	Phenanthrene-d10	17.051

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Methyl stearate	298	C19H38O2	000112-61-8	98
2			Methyl stearate	298	C19H38O2	000112-61-8	96
3			Methyl stearate	298	C19H38O2	000112-61-8	95
4			Methyl stearate	298	C19H38O2	000112-61-8	95
5			Methyl stearate	298	C19H38O2	000112-61-8	92



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN061424\
Data File : BN031996.D
Acq On : 14 Jun 2024 11:41
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Sample : P2695-17
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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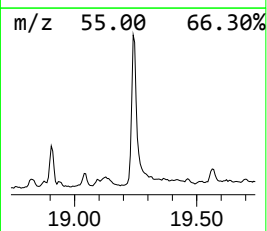
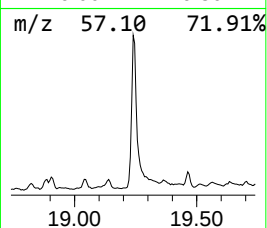
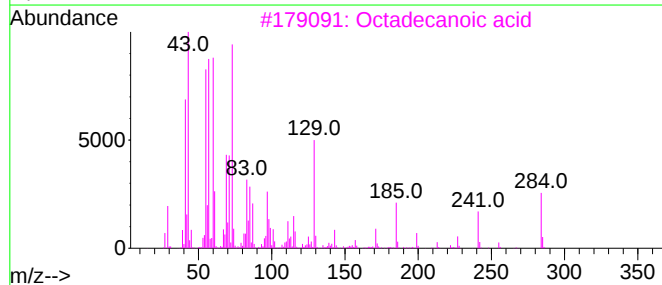
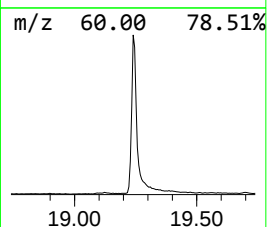
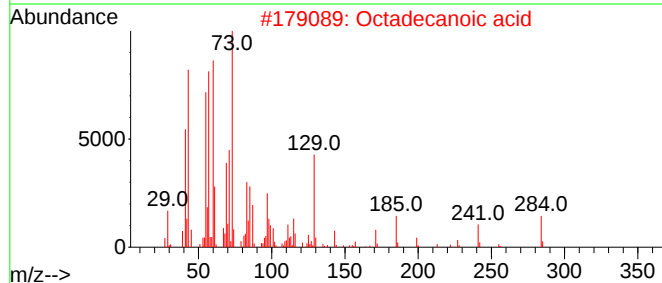
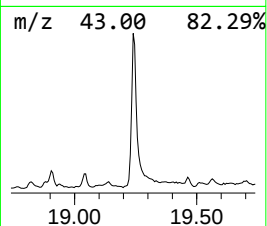
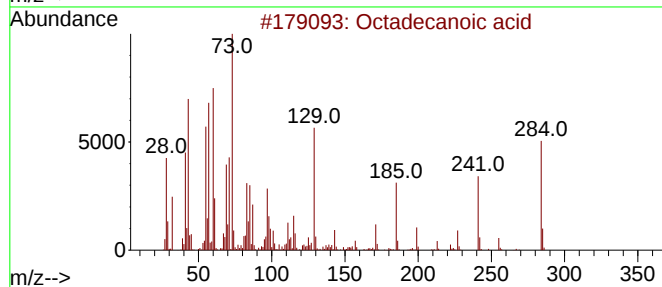
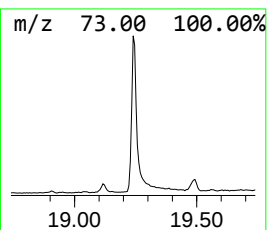
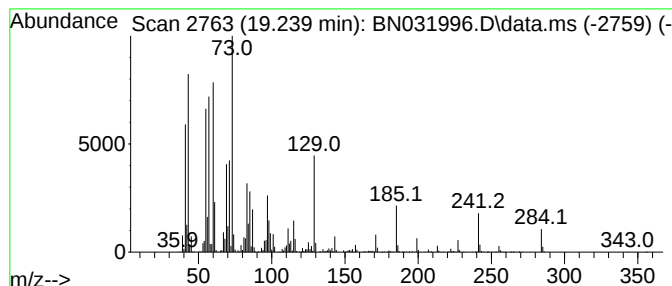
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 Octadecanoic acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.239	3.48 ng/ul	2876640	Chrysene-d12	21.292

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	95
5			Octadecanoic acid	284	C18H36O2	000057-11-4	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN061424\
Data File : BN031996.D
Acq On : 14 Jun 2024 11:41
Operator : MA/JU
Sample : P2695-17
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
A4DG5

Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN052924.MA.M
Quant Title : SVOA CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Benzene, 1,4-di...	7.540	62.1	ng/ul	5595590	1	7.658	1800630	20.0
Benzene, 1,2-di...	8.005	189.7	ng/ul	17075100	1	7.658	1800630	20.0
1-Phenoxypropan...	11.187	3.3	ng/ul	388867	2	10.440	2348420	20.0
Dodecanoic acid...	14.428	2.3	ng/ul	348014	3	14.304	2982480	20.0
Hexadecanoic ac...	17.728	3.9	ng/ul	618608	4	17.051	3140650	20.0
n-Hexadecanoic ...	17.957	23.5	ng/ul	3694820	4	17.051	3140650	20.0
9,10-Anthracene...	18.481	5.1	ng/ul	802453	4	17.051	3140650	20.0
9-Octadecenoic ...	18.904	2.7	ng/ul	430269	4	17.051	3140650	20.0
Methyl stearate	19.045	2.4	ng/ul	374286	4	17.051	3140650	20.0
Octadecanoic acid	19.239	3.5	ng/ul	2876640	5	21.292	16536900	20.0