

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN101724\
 Data File : BN034618.D
 Acq On : 17 Oct 2024 12:26
 Operator : RC/JU
 Sample : P4360-02
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A4E85

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-BN101624.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BN034618.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.964	9	13	34	rVB	2700526	3368314	14.22%	0.797%
2	3.211	51	55	59	rVB	51779	61845	0.26%	0.015%
3	6.069	530	541	547	rBV2	49715	89804	0.38%	0.021%
4	6.834	664	671	673	rVV	862713	1413934	5.97%	0.334%
5	6.857	673	675	683	rVV	1338823	1924868	8.13%	0.455%
6	7.010	693	701	709	rVV	666823	1056597	4.46%	0.250%
7	7.105	710	717	727	rVV	6006048	8799385	37.16%	2.081%
8	7.205	727	734	735	rVV	829795	1251191	5.28%	0.296%
9	7.234	735	739	749	rVV	3277085	5087716	21.48%	1.203%
10	7.669	804	813	814	rBV	824540	1240438	5.24%	0.293%
11	7.705	814	819	832	rVV	7947999	12725256	53.73%	3.010%
12	8.016	863	872	882	rBV	7989631	12812393	54.10%	3.030%
13	8.375	926	933	942	rBV	974760	1547220	6.53%	0.366%
14	8.481	942	951	965	rVB	6008173	9795865	41.36%	2.317%
15	8.740	987	995	1001	rBV	4268249	6741637	28.47%	1.595%
16	8.810	1001	1007	1010	rVV	771460	1288731	5.44%	0.305%
17	8.851	1010	1014	1021	rVV	1373124	2117845	8.94%	0.501%
18	9.387	1094	1105	1113	rBV	6995032	11771085	49.70%	2.784%
19	9.528	1121	1129	1130	rBV	501612	659679	2.79%	0.156%
20	9.557	1130	1134	1145	rVB	1174031	1951852	8.24%	0.462%
21	9.687	1150	1156	1164	rVB	301349	470537	1.99%	0.111%
22	9.875	1180	1188	1201	rBV	6545852	10346415	43.69%	2.447%
23	10.093	1213	1225	1240	rBV2	4169138	8041615	33.96%	1.902%
24	10.440	1276	1284	1291	rBV	1134099	1865220	7.88%	0.441%
25	10.569	1300	1306	1312	rBV	715568	1195288	5.05%	0.283%
26	10.793	1332	1344	1354	rBV	5525392	8756057	36.97%	2.071%
27	11.722	1488	1502	1510	rBV	1705860	2668580	11.27%	0.631%
28	11.998	1544	1549	1553	rBV	66173	96000	0.41%	0.023%
29	12.110	1560	1568	1576	rBV	3890774	5991660	25.30%	1.417%
30	12.540	1636	1641	1647	rBV	106964	154869	0.65%	0.037%
31	12.722	1664	1672	1678	rBV	6370584	10044343	42.41%	2.376%
32	12.787	1678	1683	1694	rVB	1588103	2377358	10.04%	0.562%
33	13.163	1742	1747	1750	rVV	55574	78705	0.33%	0.019%
34	13.210	1750	1755	1760	rVB3	32682	61729	0.26%	0.015%
35	13.728	1836	1843	1850	rBV	1363306	2011036	8.49%	0.476%
36	13.881	1861	1869	1877	rBV	7285919	10846157	45.80%	2.565%

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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_N\Methods\SFAM-EPA-BN101624.MA.M
 Title : SVOA CALIBRATION

37	13.987	1880	1887	1890	rVV2	1856460	3023115	12.77%	0.715%
38	14.016	1890	1892	1899	rVB	1493607	1813857	7.66%	0.429%
39	14.298	1934	1940	1945	rBV	1460598	2099312	8.86%	0.497%
40	14.369	1945	1952	1958	rVB	8872714	13211904	55.79%	3.125%
41	14.510	1963	1976	1990	rBV2	2647027	4707018	19.88%	1.113%
42	14.616	1990	1994	1996	rVV	106687	145896	0.62%	0.035%
43	14.669	1996	2003	2013	rVV	5111553	7337449	30.98%	1.735%
44	14.922	2041	2046	2052	rBV2	79974	121676	0.51%	0.029%
45	15.151	2078	2085	2090	rBV6	30146	67333	0.28%	0.016%
46	15.298	2103	2110	2114	rBV	2157327	3161767	13.35%	0.748%
47	15.357	2114	2120	2125	rVV	14195381	20804182	87.85%	4.921%
48	15.404	2125	2128	2138	rVB	395583	494630	2.09%	0.117%
49	15.604	2156	2162	2166	rVV	200484	276331	1.17%	0.065%
50	15.663	2166	2172	2180	rVB	904932	1217555	5.14%	0.288%
51	15.945	2215	2220	2225	rBV5	38420	62386	0.26%	0.015%
52	16.010	2225	2231	2236	rVV3	86510	150285	0.63%	0.036%
53	16.245	2262	2271	2279	rBV2	2287854	3273418	13.82%	0.774%
54	16.357	2282	2290	2296	rVV	9539786	14384768	60.74%	3.402%
55	16.698	2339	2348	2355	rBV	6695111	10247243	43.27%	2.424%
56	16.857	2372	2375	2381	rVB	65022	87676	0.37%	0.021%
57	17.045	2400	2407	2410	rBV	1577639	2382081	10.06%	0.563%
58	17.086	2410	2414	2419	rVV	3834627	5298256	22.37%	1.253%
59	17.139	2419	2423	2426	rVV2	2286849	3321342	14.02%	0.786%
60	17.175	2426	2429	2439	rVB	3961303	5085812	21.47%	1.203%
61	17.663	2508	2512	2515	rBV4	58245	88095	0.37%	0.021%
62	17.692	2515	2517	2522	rVB2	52036	70209	0.30%	0.017%
63	17.757	2523	2528	2533	rBV	185904	226090	0.95%	0.053%
64	17.975	2560	2565	2570	rBV	327916	467907	1.98%	0.111%
65	18.045	2571	2577	2582	rVV	11210276	14559065	61.48%	3.443%
66	18.086	2582	2584	2592	rVV4	89208	145804	0.62%	0.034%
67	18.269	2610	2615	2618	rBV3	68741	113482	0.48%	0.027%
68	18.327	2622	2625	2635	rVB6	50411	118850	0.50%	0.028%
69	18.410	2635	2639	2642	rBV2	60103	90367	0.38%	0.021%
70	18.451	2642	2646	2652	rVB	409426	536421	2.27%	0.127%
71	18.810	2703	2707	2716	rVB3	118989	252948	1.07%	0.060%
72	18.886	2717	2720	2725	rBV3	40889	82721	0.35%	0.020%
73	18.933	2725	2728	2732	rVB3	73034	83717	0.35%	0.020%
74	19.110	2744	2758	2763	rBV	11955765	17862838	75.43%	4.225%
75	19.257	2779	2783	2792	rBV6	83817	158111	0.67%	0.037%
76	19.339	2793	2797	2803	rVB	123033	162059	0.68%	0.038%

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77	19.433	2807	2813	2815	rBV	2767631	3838594	16.21%	0.908%
78	19.463	2815	2818	2824	rVB	6510587	8439510	35.64%	1.996%
79	19.592	2838	2840	2845	rVB4	68602	87521	0.37%	0.021%
80	19.998	2906	2909	2913	rBV5	53277	79748	0.34%	0.019%
81	20.151	2932	2935	2937	rBV	131516	145133	0.61%	0.034%
82	20.174	2937	2939	2946	rVB	108484	135696	0.57%	0.032%
83	20.392	2972	2976	2981	rBV	3563377	3831224	16.18%	0.906%
84	20.998	3075	3079	3093	rBV2	571638	892766	3.77%	0.211%
85	21.157	3103	3106	3112	rVB	219311	250985	1.06%	0.059%
86	21.227	3113	3118	3121	rBV	14829335	22169304	93.61%	5.243%
87	21.257	3121	3123	3129	rVV2	10691297	14924573	63.02%	3.530%
88	21.321	3129	3134	3140	rVB	15623689	23682703	100.00%	5.601%
89	22.092	3261	3265	3277	rBV2	166464	295166	1.25%	0.070%
90	22.274	3292	3296	3305	rVB2	165941	269026	1.14%	0.064%
91	23.268	3454	3465	3484	rVB	10248239	23453055	99.03%	5.547%
92	23.463	3493	3498	3502	rBV	153959	338036	1.43%	0.080%
93	23.962	3574	3583	3588	rBV	1630868	3877520	16.37%	0.917%
94	24.027	3588	3594	3606	rVV	2454007	5530805	23.35%	1.308%
95	24.157	3608	3616	3632	rVB	1143397	2764526	11.67%	0.654%
96	25.357	3814	3820	3827	rBV	136000	329280	1.39%	0.078%
97	25.445	3828	3835	3850	rVB2	199149	611637	2.58%	0.145%
98	27.398	4154	4167	4169	rBV2	791039	2381654	10.06%	0.563%
99	27.480	4171	4181	4197	rVB	3296180	11958934	50.50%	2.829%
100	28.398	4323	4337	4355	rVB	963287	4007273	16.92%	0.948%

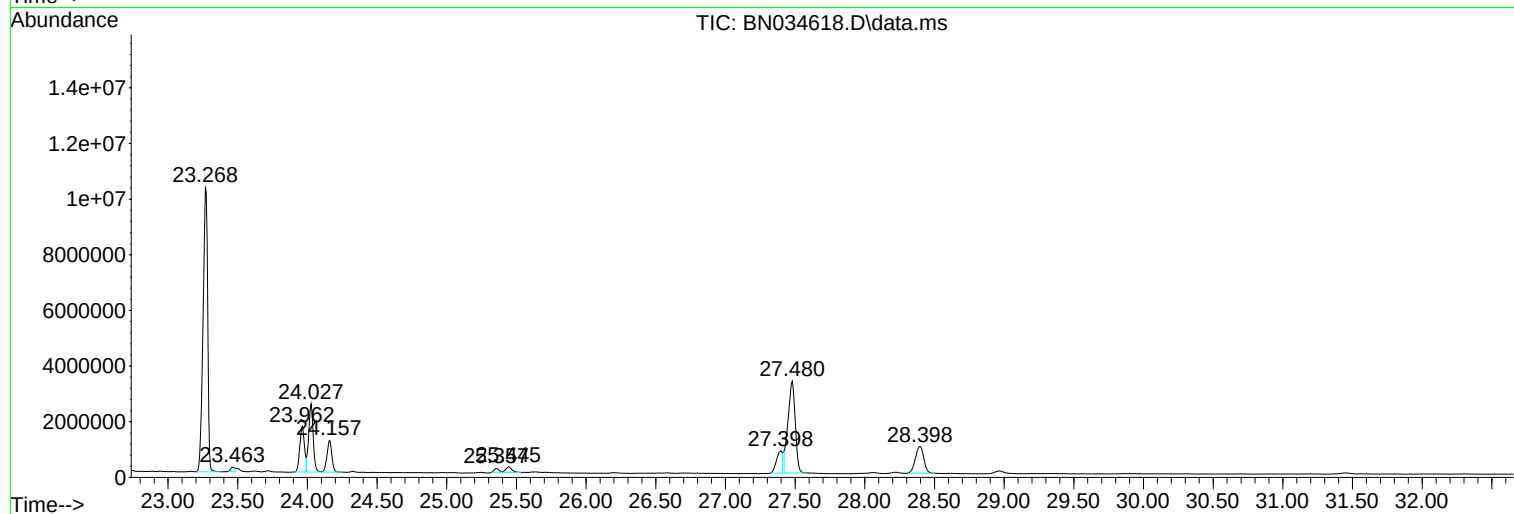
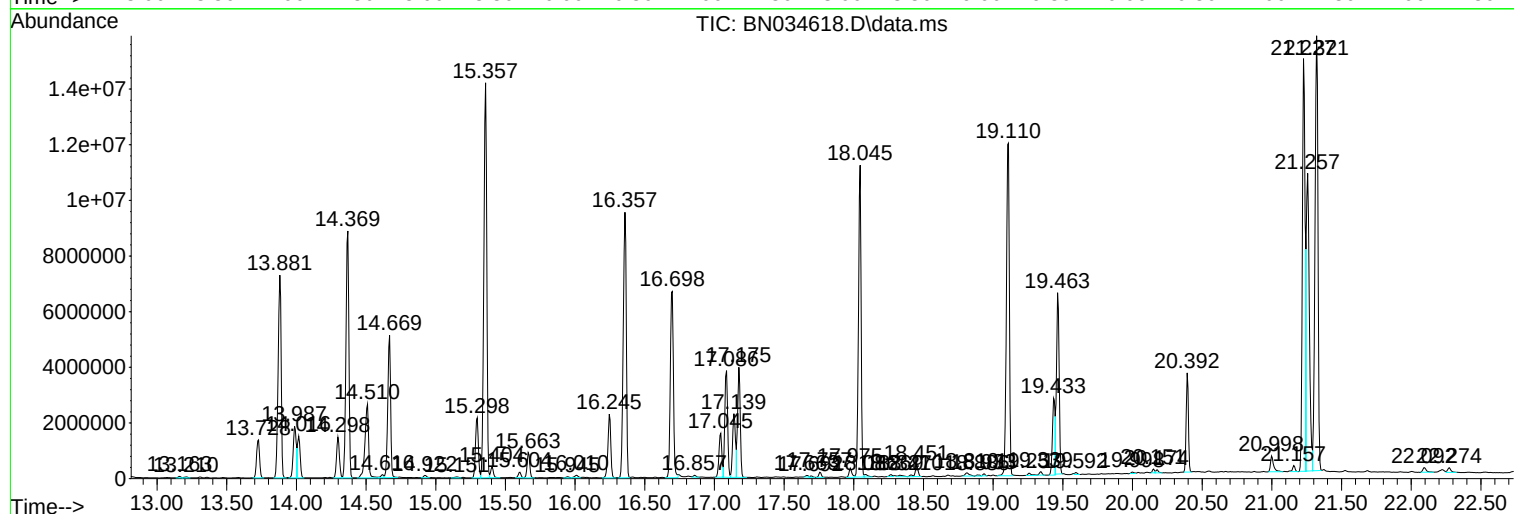
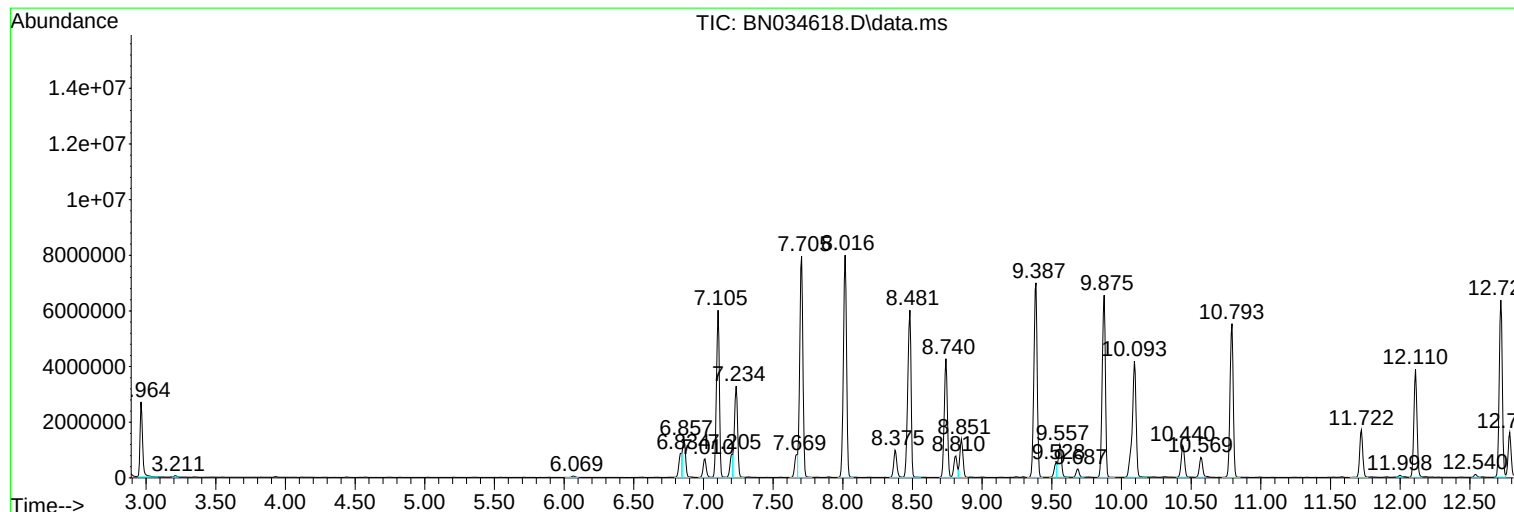
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Instrument :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN101624.MA.M
Quant Title : SVOA CALIBRATION

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



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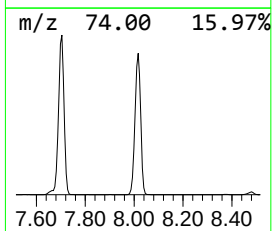
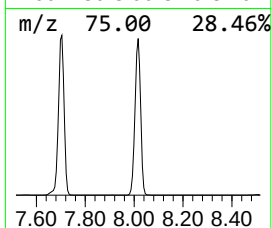
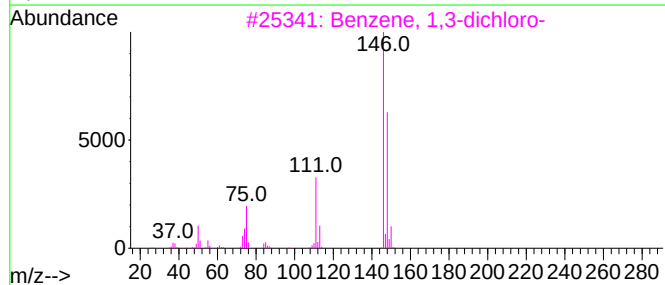
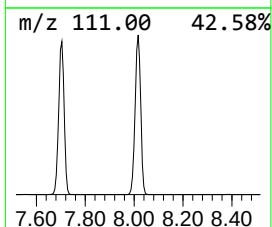
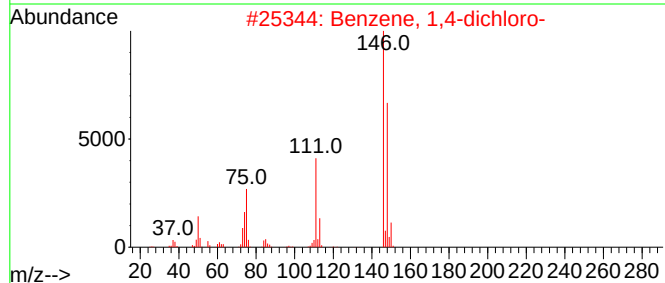
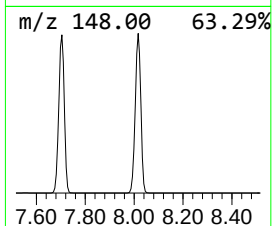
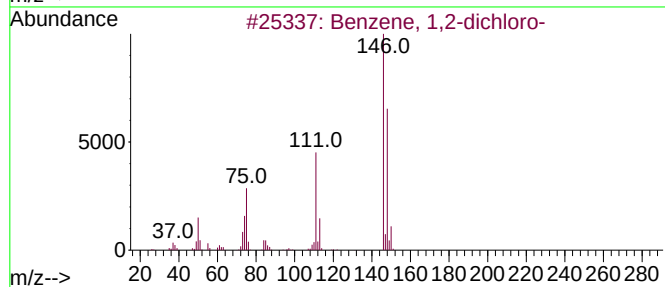
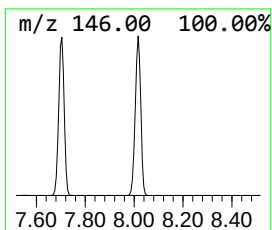
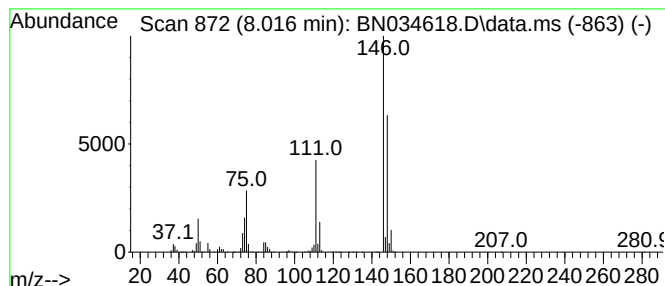
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN101624.MA.M
Quant Title : SVOA CALIBRATION

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.
8.016	206.58 ng/ul	12812400	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	97
4			Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	97
5			Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	97



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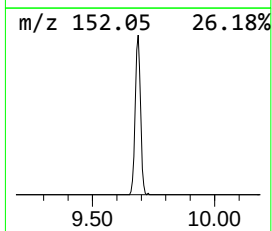
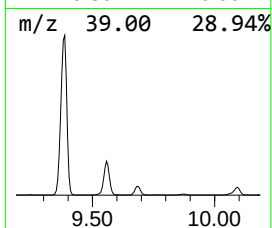
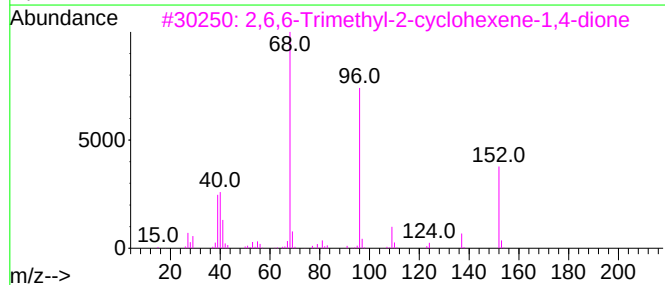
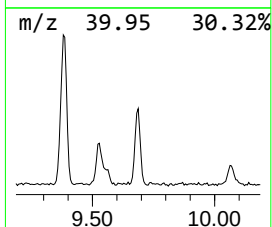
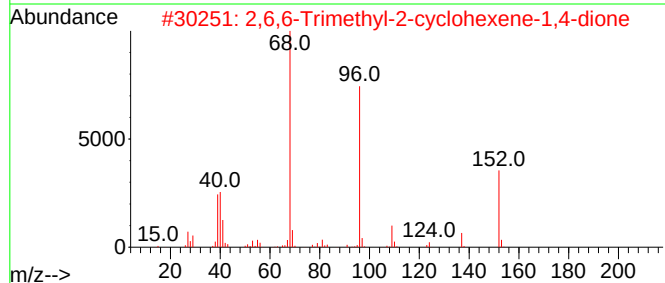
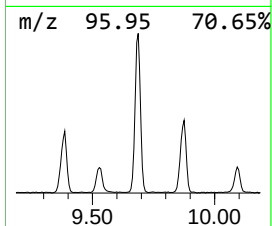
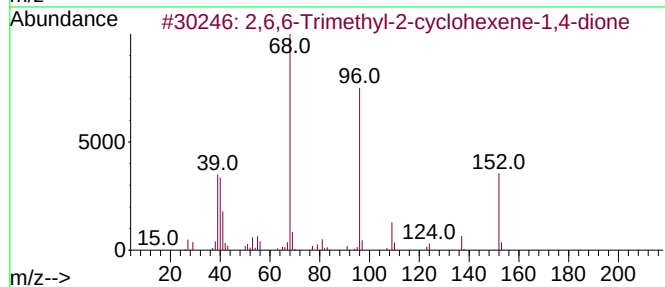
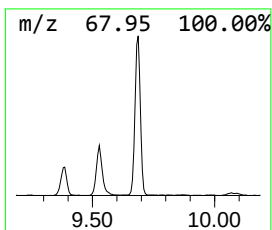
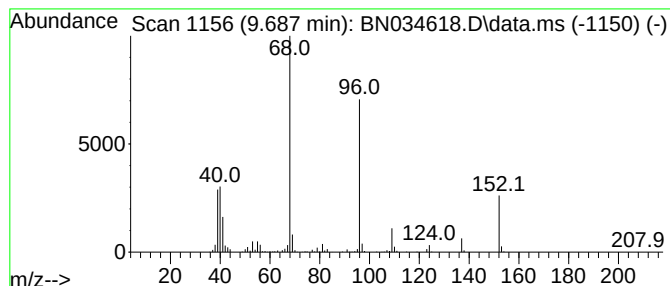
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN101624.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 2,6,6-Trimethyl-2-cyclohexe... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.687	5.05 ng/ul	470537	Naphthalene-d8	10.440

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	53		



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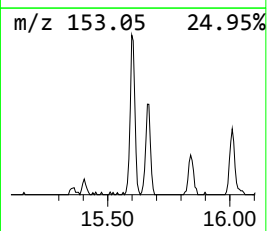
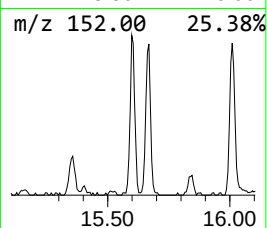
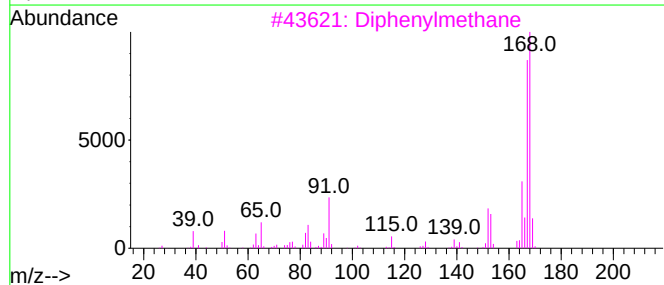
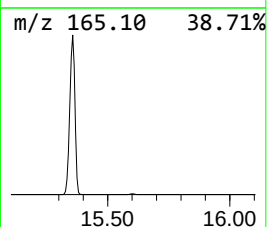
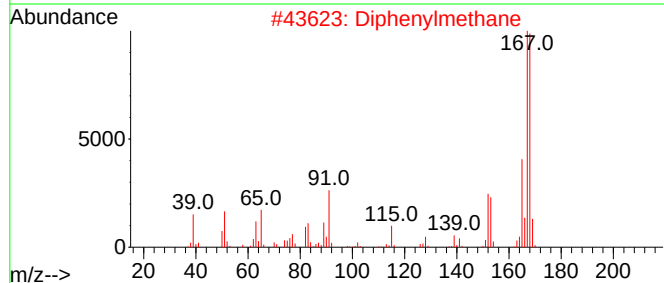
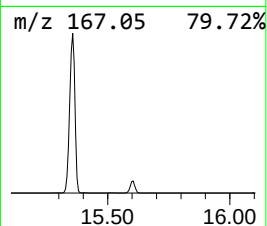
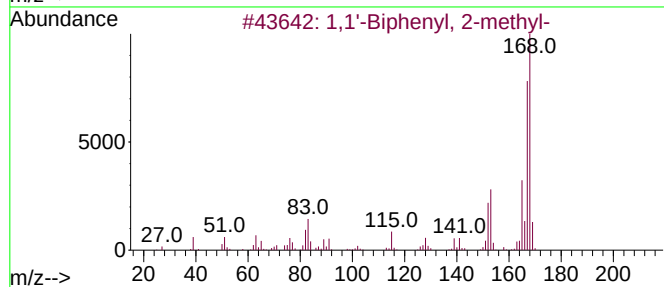
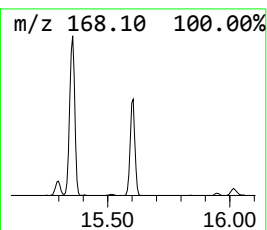
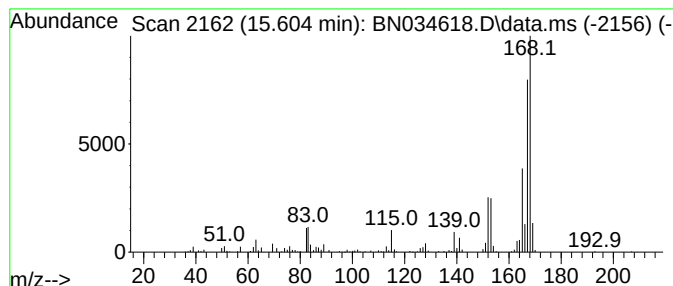
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 1,1'-Biphenyl, 2-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.604	2.63 ng/ul	276331	Acenaphthene-d10	14.298

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,1'-Biphenyl, 2-methyl-	168	C13H12	000643-58-3	96
2			Diphenylmethane	168	C13H12	000101-81-5	93
3			Diphenylmethane	168	C13H12	000101-81-5	91
4			Diphenylmethane	168	C13H12	000101-81-5	91
5			1,1'-Biphenyl, 2-methyl-	168	C13H12	000643-58-3	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN101724\
Data File : BN034618.D
Acq On : 17 Oct 2024 12:26
Operator : RC/JU
Sample : P4360-02
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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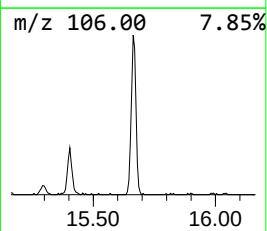
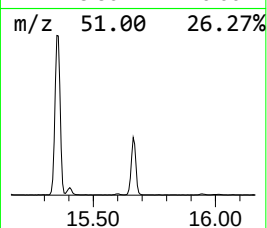
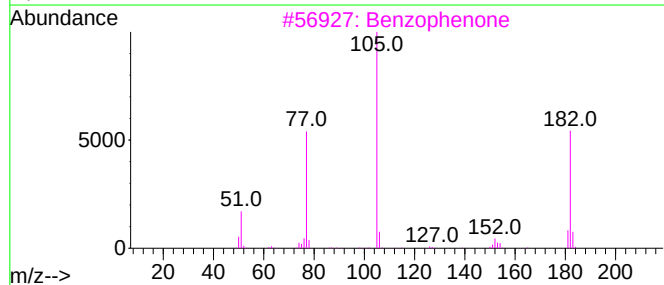
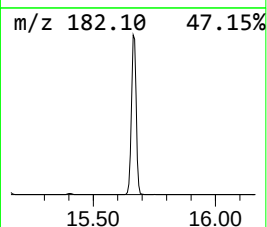
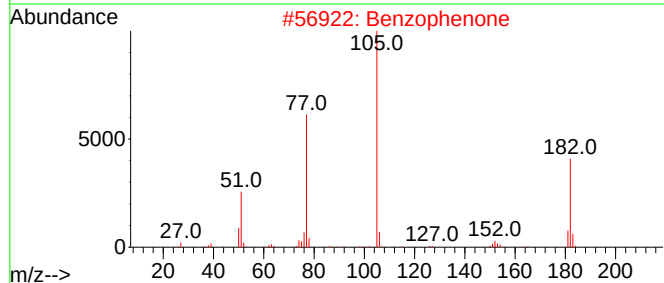
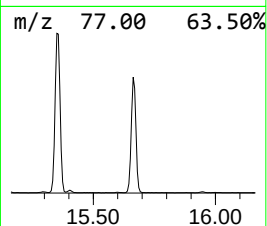
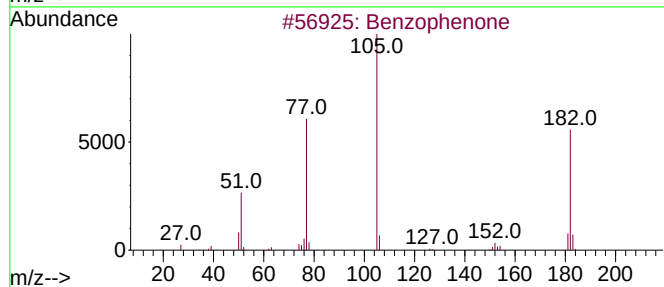
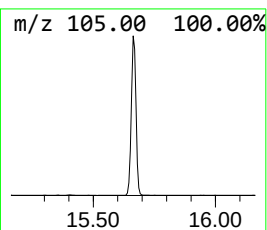
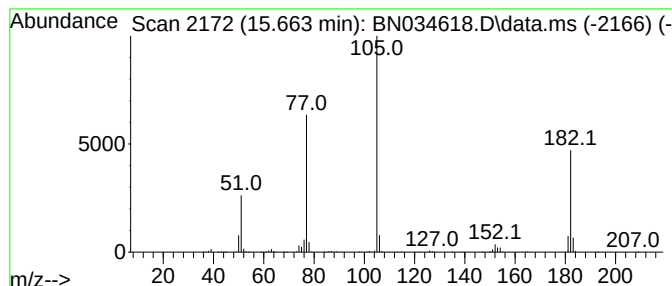
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Benzophenone Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.663	11.60 ng/ul	1217560	Acenaphthene-d10	14.298

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	96
2			Benzophenone	182	C13H10O	000119-61-9	96
3			Benzophenone	182	C13H10O	000119-61-9	95
4			Benzophenone	182	C13H10O	000119-61-9	94
5			1,2,4-Trioxolane, 3,3,5-triphenyl-	304	C20H16O3	023246-12-0	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN101724\
Data File : BN034618.D
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Sample : P4360-02
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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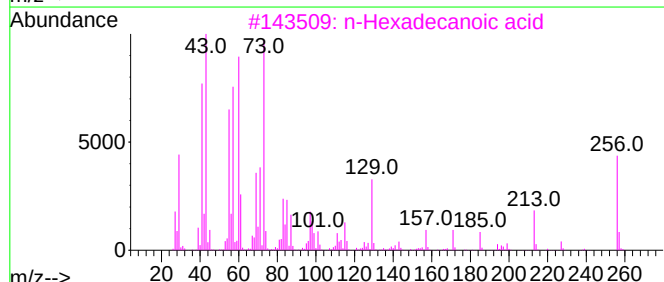
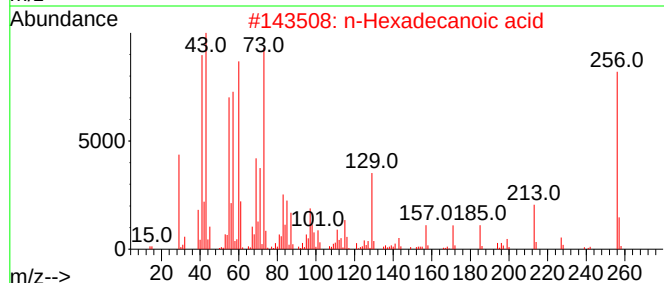
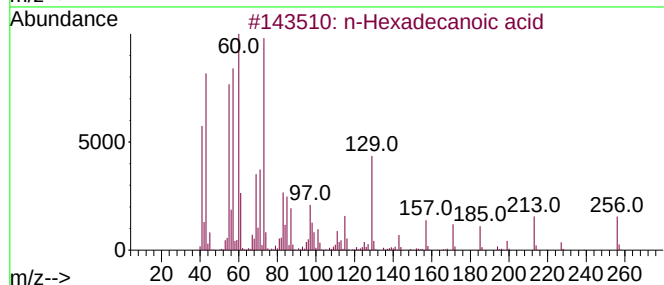
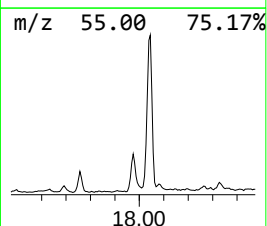
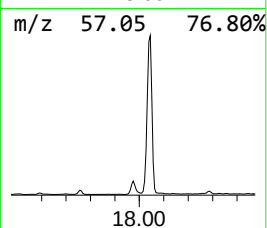
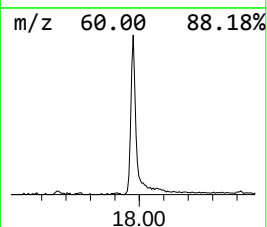
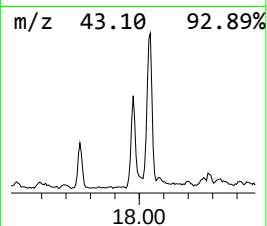
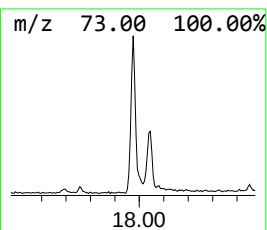
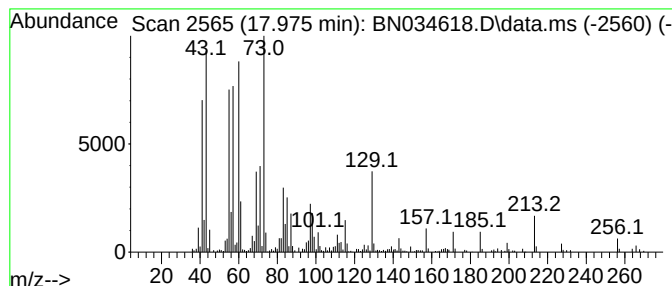
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 n-Hexadecanoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.975	3.93 ng/ul	467907	Phenanthrene-d10	17.045

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
2			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
3			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
5			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN101724\
Data File : BN034618.D
Acq On : 17 Oct 2024 12:26
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Sample : P4360-02
Misc :
ALS Vial : 6 Sample Multiplier: 1

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ClientSampleId :
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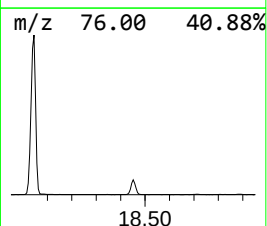
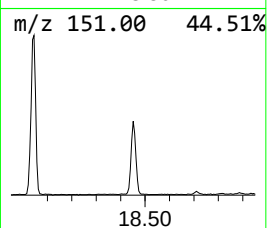
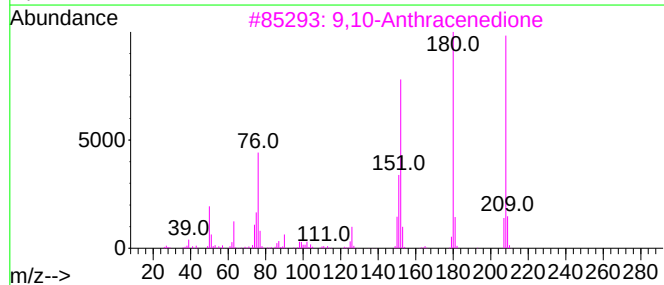
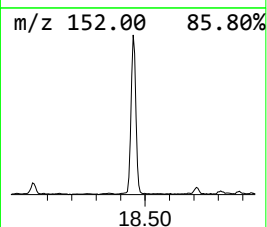
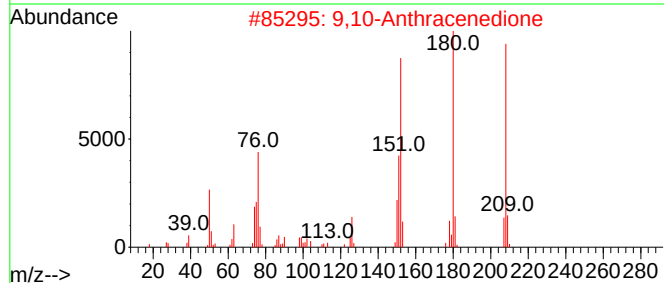
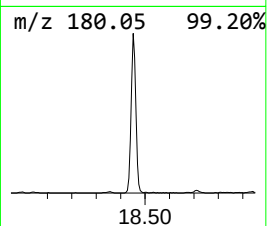
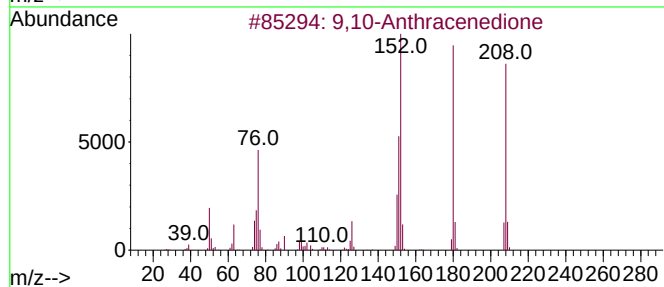
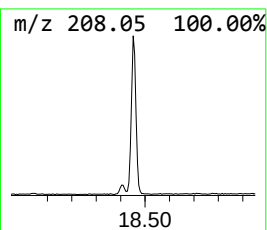
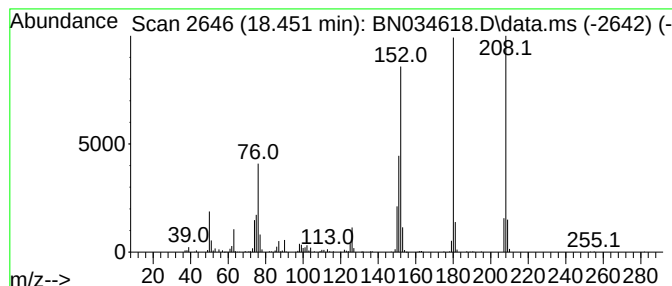
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 9,10-Anthracenedione Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.451	4.50 ng/ul	536421	Phenanthrene-d10	17.045

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	96
5	9H-Fluoren-9-one			180	C13H8O	000486-25-9	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN101724\
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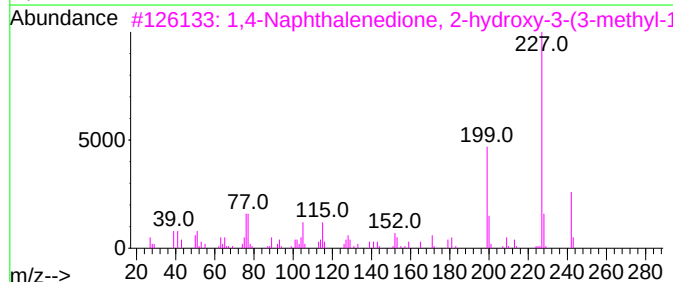
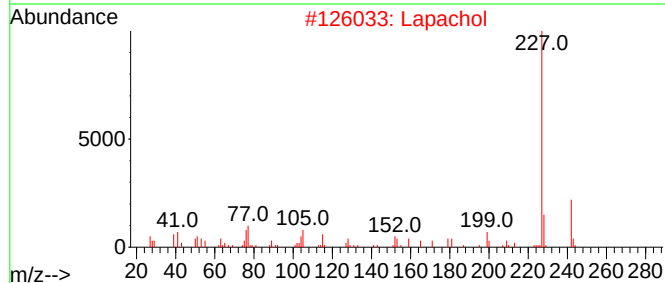
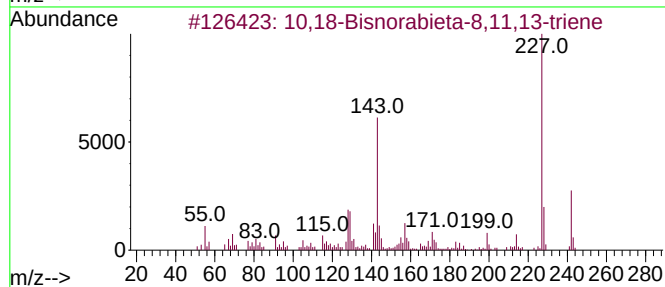
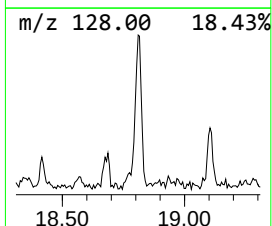
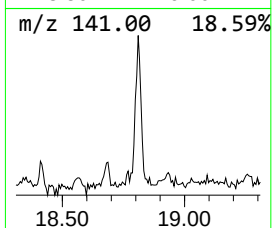
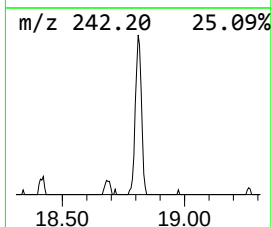
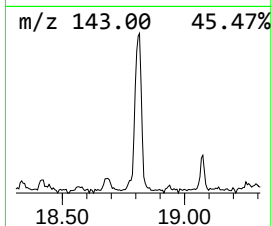
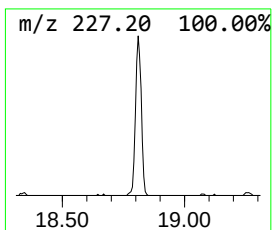
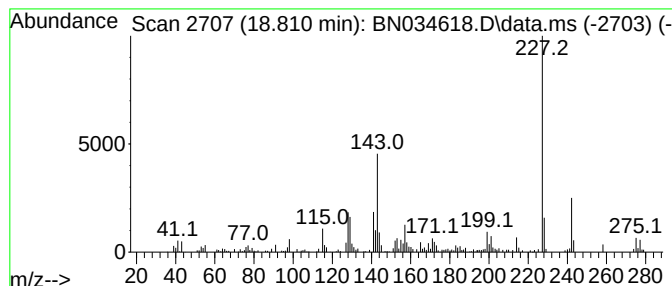
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 10,18-Bisnorabieta-8,11,13-... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD		R.T.	
18.810	2.12 ng/ul	252948	Phenanthrene-d10		17.045	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	10,18-Bisnorabieta-8,11,13-triene		242	C18H26	032624-67-2	99
2	Lapachol		242	C15H14O3	000084-79-7	64
3	1,4-Naphthalenedione, 2-hydroxy-...		242	C15H14O3	004042-39-1	53
4	s-Indacene-1,7-dione, 2,3,5,6-te...		242	C16H18O2	055591-17-8	53
5	5-Methoxy-2-methyl-4-oxo-1,2,3,4...		242	C14H14N2O2	1000227-06-0	52



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN101724\
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ALS Vial : 6 Sample Multiplier: 1

Instrument :
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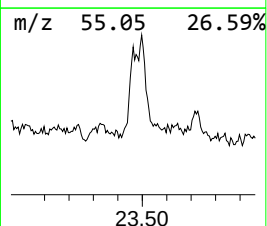
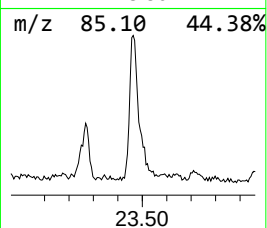
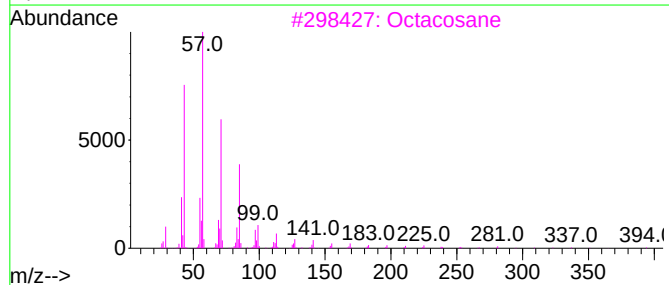
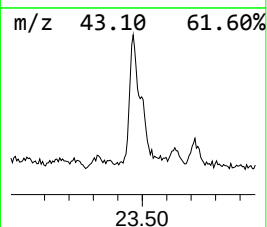
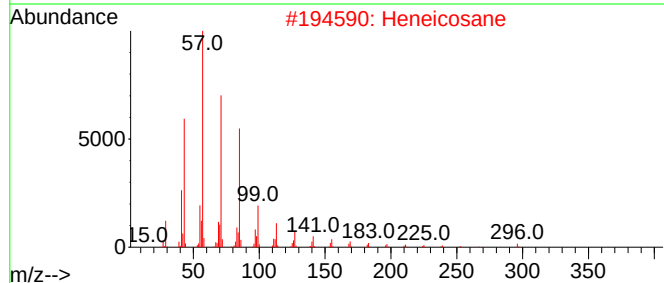
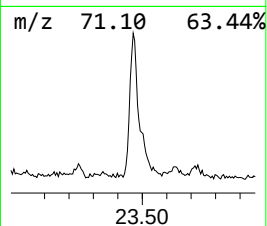
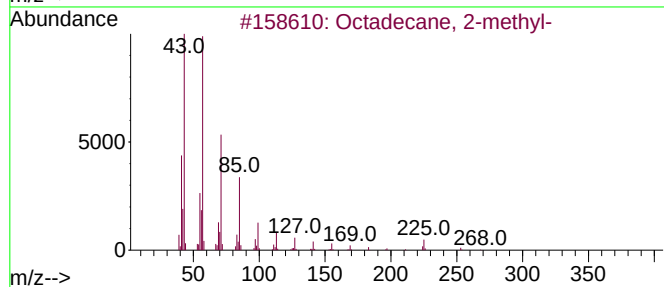
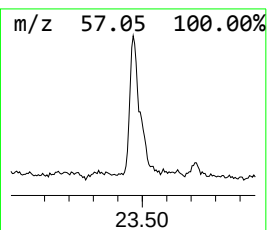
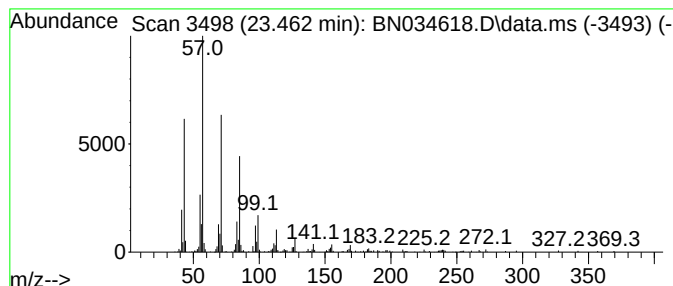
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.462	2.45 ng/ul	338036	Perylene-d12	24.157

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Octadecane, 2-methyl-	268	C19H40	001560-88-9	93
2		Heneicosane	296	C21H44	000629-94-7	91
3		Octacosane	394	C28H58	000630-02-4	91
4		Tetracosane	338	C24H50	000646-31-1	90
5		Tetradecane, 2,6,10-trimethyl-	240	C17H36	014905-56-7	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN101724\
Data File : BN034618.D
Acq On : 17 Oct 2024 12:26
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Misc :
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Instrument :
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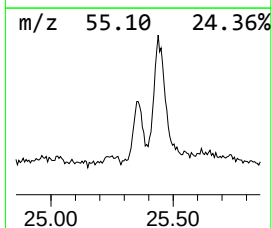
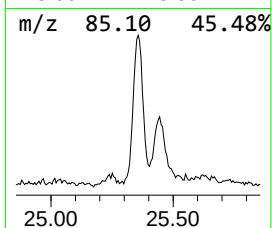
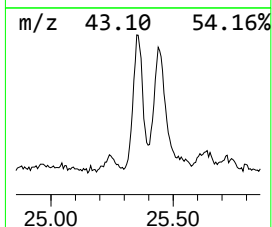
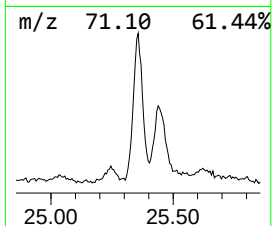
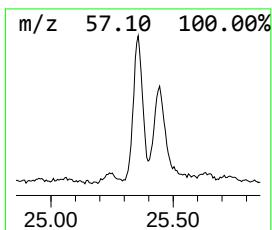
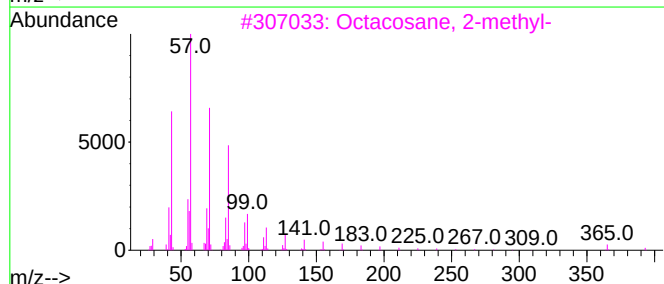
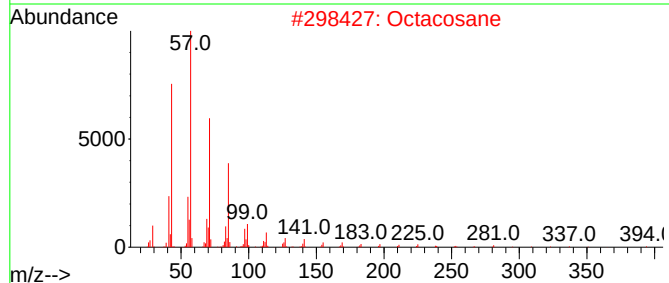
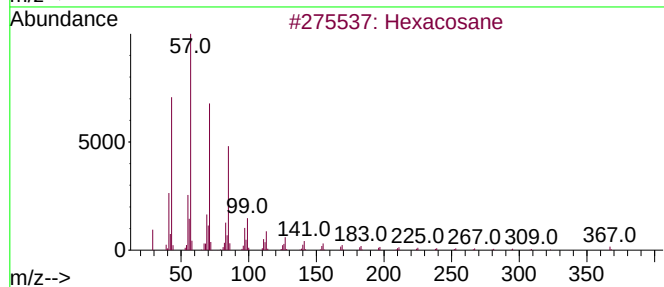
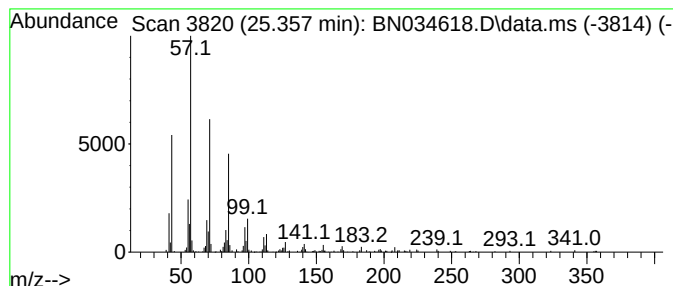
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.357	2.38 ng/ul	329280	Perylene-d12	24.157

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Hexacosane	366	C26H54	000630-01-3	91
2		Octacosane	394	C28H58	000630-02-4	90
3		Octacosane, 2-methyl-	408	C29H60	001560-98-1	87
4		Tridecane, 2-methyl-	198	C14H30	001560-96-9	87
5		Sulfurous acid, butyl dodecyl ester	306	C16H34O3S	1000309-17-9	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN101724\
Data File : BN034618.D
Acq On : 17 Oct 2024 12:26
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Misc :
ALS Vial : 6 Sample Multiplier: 1

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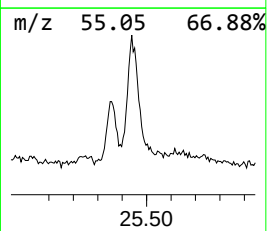
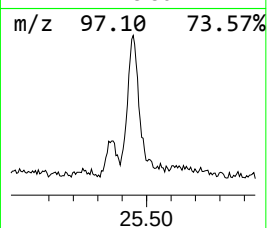
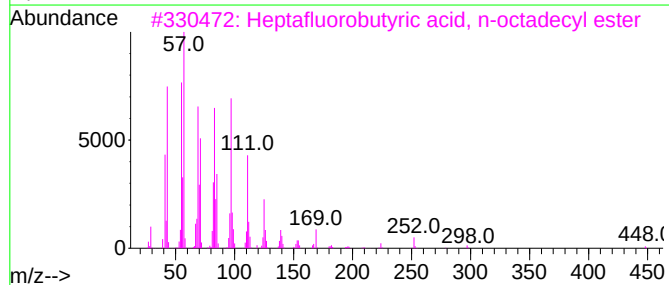
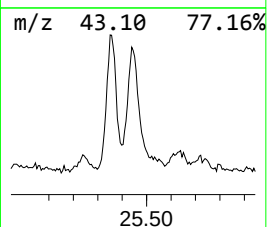
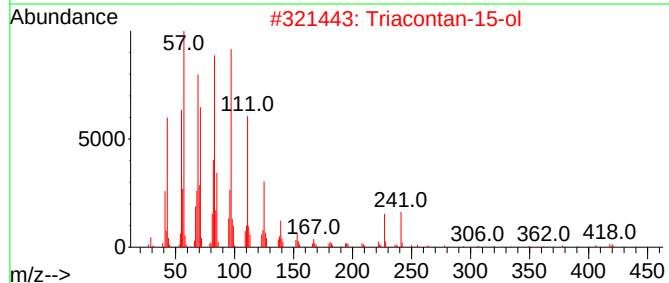
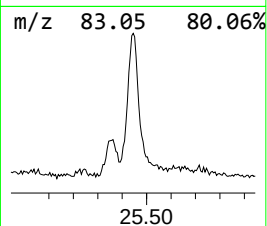
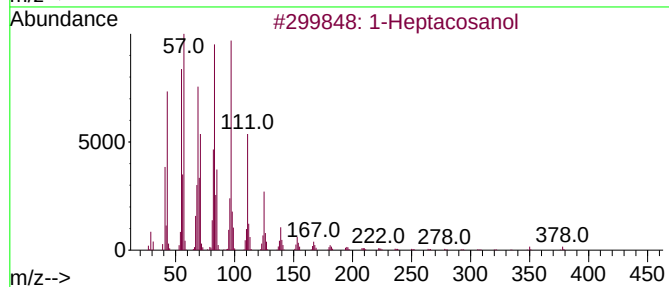
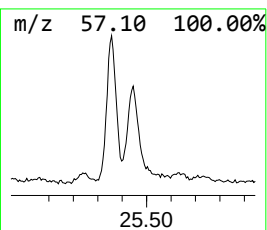
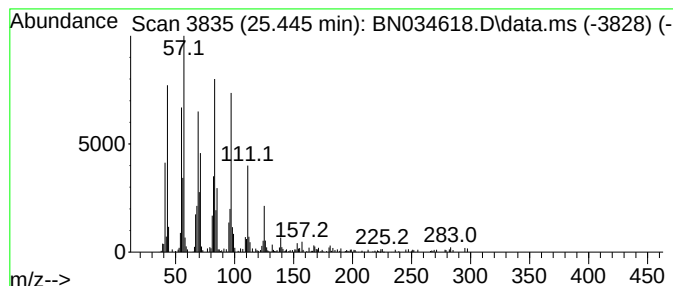
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 1-Heptacosanol Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.445	4.42 ng/ul	611637	Perylene-d12	24.157

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Heptacosanol	396	C27H56O	002004-39-9	95
2			Triacontan-15-ol	438	C30H62O	031849-26-0	91
3			Heptafluorobutyric acid, n-octadec...	466	C22H37F7O2	000400-57-7	91
4			17-Pentatriacontene	491	C35H70	006971-40-0	90
5			1-Tetracosene	336	C24H48	010192-32-2	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN101724\
Data File : BN034618.D
Acq On : 17 Oct 2024 12:26
Operator : RC/JU
Sample : P4360-02
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
A4E85

Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN101624.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
Benzene, 1,2-di...	8.016	206.6	ng/ul	12812400	1	7.669	1240440	20.0
2,6,6-Trimethyl...	9.687	5.0	ng/ul	470537	2	10.440	1865220	20.0
1,1'-Biphenyl, ...	15.604	2.6	ng/ul	276331	3	14.298	2099310	20.0
Benzophenone	15.663	11.6	ng/ul	1217560	3	14.298	2099310	20.0
n-Hexadecanoic ...	17.975	3.9	ng/ul	467907	4	17.045	2382080	20.0
9,10-Anthracene...	18.451	4.5	ng/ul	536421	4	17.045	2382080	20.0
10,18-Bisnorabi...	18.810	2.1	ng/ul	252948	4	17.045	2382080	20.0
(DEL) Alkane: S...	23.462	2.5	ng/ul	338036	6	24.157	2764530	20.0
(DEL) Alkane: S...	25.357	2.4	ng/ul	329280	6	24.157	2764530	20.0
1-Heptacosanol	25.445	4.4	ng/ul	611637	6	24.157	2764530	20.0