

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062224\
 Data File : BN032177.D
 Acq On : 22 Jun 2024 21:50
 Operator : MA/JU
 Sample : P2899-18
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0AB9

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: BN032177.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.163	26	30	43	rVB	70144	110506	1.75%	0.135%
2	3.434	72	76	84	rVB	75974	107152	1.70%	0.131%
3	3.581	98	101	111	rVV	103263	145073	2.30%	0.178%
4	3.669	112	116	123	rVB	86332	110362	1.75%	0.135%
5	3.875	147	151	157	rVB	34408	48242	0.77%	0.059%
6	4.375	232	236	240	rVV	16525	29583	0.47%	0.036%
7	4.422	240	244	249	rVV	57233	86645	1.38%	0.106%
8	4.475	249	253	260	rVB	25840	41787	0.66%	0.051%
9	4.769	298	303	313	rBV	327398	486516	7.72%	0.596%
10	5.010	340	344	351	rVB3	17305	28868	0.46%	0.035%
11	5.293	388	392	399	rBV2	15545	30625	0.49%	0.038%
12	5.657	447	454	459	rVV4	29272	52011	0.83%	0.064%
13	5.734	459	467	477	rVB2	47040	90086	1.43%	0.110%
14	6.157	528	539	550	rBV6	18500	54811	0.87%	0.067%
15	6.316	560	566	576	rVB3	60816	113056	1.80%	0.138%
16	6.663	620	625	631	rVB3	26988	49956	0.79%	0.061%
17	6.816	639	651	663	rVB	1002101	1777563	28.22%	2.178%
18	6.975	673	678	689	rVB	881494	1407120	22.34%	1.724%
19	7.063	689	693	704	rVB2	32256	61484	0.98%	0.075%
20	7.169	704	711	720	rBV	1167802	1908644	30.31%	2.338%
21	7.245	720	724	727	rBV	35459	49594	0.79%	0.061%
22	7.634	776	790	797	rBV	1118066	1833069	29.11%	2.246%
23	7.704	797	802	806	rVV2	27105	57853	0.92%	0.071%
24	7.898	829	835	840	rBV2	25778	44155	0.70%	0.054%
25	8.345	905	911	922	rVV	1096313	1859076	29.52%	2.277%
26	8.457	924	930	938	rVV	202898	356748	5.66%	0.437%
27	8.792	979	987	993	rBV	1033246	1749772	27.78%	2.144%
28	9.357	1077	1083	1089	rVB	98014	155912	2.48%	0.191%
29	9.504	1099	1108	1126	rBV	853404	1570958	24.94%	1.925%
30	9.692	1132	1140	1148	rBV	62173	177260	2.81%	0.217%
31	9.863	1165	1169	1176	rVB2	25916	48244	0.77%	0.059%
32	10.045	1193	1200	1221	rVV	1291172	2368187	37.60%	2.901%
33	10.192	1221	1225	1234	rVB	13162	30248	0.48%	0.037%
34	10.416	1252	1263	1268	rBV	1407549	2484511	39.45%	3.044%
35	10.463	1268	1271	1276	rVV	45707	71482	1.14%	0.088%
36	10.563	1282	1288	1307	rVB	174222	380398	6.04%	0.466%

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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-BN052924.MA.M
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37	10.986	1345	1360	1377	rBV	153388	713629	11.33%	0.874%
38	11.163	1384	1390	1395	rBV	163651	295297	4.69%	0.362%
39	11.216	1396	1399	1410	rVB3	39569	77678	1.23%	0.095%
40	11.422	1430	1434	1444	rVB3	28242	65163	1.03%	0.080%
41	11.828	1499	1503	1514	rVB2	46064	80613	1.28%	0.099%
42	12.004	1526	1533	1540	rBV2	37153	67306	1.07%	0.082%
43	12.080	1541	1546	1551	rVB	74544	112559	1.79%	0.138%
44	12.304	1579	1584	1590	rVB	60780	99167	1.57%	0.121%
45	12.404	1595	1601	1608	rBV	41807	78869	1.25%	0.097%
46	12.551	1621	1626	1632	rVV6	15659	29885	0.47%	0.037%
47	12.792	1665	1667	1673	rVB	19936	29390	0.47%	0.036%
48	13.092	1713	1718	1723	rBV2	54883	95207	1.51%	0.117%
49	13.139	1723	1726	1733	rVV3	28947	50151	0.80%	0.061%
50	13.210	1733	1738	1744	rVV3	24649	55195	0.88%	0.068%
51	13.269	1744	1748	1753	rVV3	34825	62511	0.99%	0.077%
52	13.533	1788	1793	1798	rBV	186659	283151	4.50%	0.347%
53	13.592	1798	1803	1808	rVB3	78524	137258	2.18%	0.168%
54	13.651	1808	1813	1814	rBV	41804	55236	0.88%	0.068%
55	13.692	1814	1820	1827	rVV	2189198	3242785	51.49%	3.973%
56	13.751	1827	1830	1834	rVV3	39153	59528	0.95%	0.073%
57	13.792	1834	1837	1840	rVV5	20811	29981	0.48%	0.037%
58	13.833	1840	1844	1847	rVV	32164	42449	0.67%	0.052%
59	13.969	1856	1867	1880	rBV	2759179	4214410	66.92%	5.163%
60	14.110	1885	1891	1895	rBV6	25199	52895	0.84%	0.065%
61	14.174	1895	1902	1906	rVB3	62295	114121	1.81%	0.140%
62	14.280	1913	1920	1926	rBV	2078747	3123066	49.59%	3.826%
63	14.339	1926	1930	1933	rBV3	35555	50840	0.81%	0.062%
64	14.516	1950	1960	1978	rBV	943747	1910380	30.33%	2.340%
65	14.663	1980	1985	1990	rBV3	88410	177149	2.81%	0.217%
66	14.892	2022	2024	2030	rVB4	17705	28781	0.46%	0.035%
67	14.951	2030	2034	2038	rVB2	24517	31630	0.50%	0.039%
68	15.074	2047	2055	2058	rBV4	40928	88328	1.40%	0.108%
69	15.127	2061	2064	2068	rVB	50114	60290	0.96%	0.074%
70	15.274	2080	2089	2095	rBV	3368445	4929824	78.28%	6.039%
71	15.410	2107	2112	2118	rBV	825528	1191845	18.92%	1.460%
72	15.627	2144	2149	2158	rBV2	39288	80917	1.28%	0.099%
73	15.768	2171	2173	2181	rVB2	27877	51031	0.81%	0.063%
74	15.851	2182	2187	2193	rVB9	33062	65624	1.04%	0.080%
75	15.998	2207	2212	2213	rBV2	98146	133168	2.11%	0.163%
76	16.151	2235	2238	2241	rBV5	25849	41422	0.66%	0.051%

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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

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 Title : SVOA CALIBRATION

77	16.298	2259	2263	2266	rBV5	25336	45541	0.72%	0.056%
78	16.439	2283	2287	2294	rBV6	60238	110423	1.75%	0.135%
79	16.686	2326	2329	2335	rBV6	30672	59503	0.94%	0.073%
80	16.786	2343	2346	2349	rBV	62799	65843	1.05%	0.081%
81	16.933	2368	2371	2376	rBV2	65040	96088	1.53%	0.118%
82	17.027	2381	2387	2392	rVV	2355109	3511892	55.76%	4.302%
83	17.068	2392	2394	2398	rVV	79347	106211	1.69%	0.130%
84	17.127	2398	2404	2414	rVB	3463972	5079144	80.65%	6.222%
85	17.210	2414	2418	2424	rBV6	46445	97872	1.55%	0.120%
86	17.315	2433	2436	2440	rBV6	28613	46899	0.74%	0.057%
87	17.527	2468	2472	2477	rBV2	75113	104304	1.66%	0.128%
88	17.874	2527	2531	2534	rBV3	71531	115006	1.83%	0.141%
89	17.933	2536	2541	2551	rVV	1307598	2009177	31.90%	2.461%
90	18.198	2582	2586	2588	rBV	181932	240786	3.82%	0.295%
91	18.545	2641	2645	2650	rVB	638304	852265	13.53%	1.044%
92	18.604	2653	2655	2660	rBV5	76780	122175	1.94%	0.150%
93	19.221	2756	2760	2764	rBV	1309165	1905698	30.26%	2.335%
94	19.427	2791	2795	2802	rVB	3840132	6297958	100.00%	7.715%
95	20.309	2941	2945	2949	rBV	2745971	3587002	56.96%	4.394%
96	20.356	2949	2953	2959	rVB	1425779	2252114	35.76%	2.759%
97	20.927	3046	3050	3052	rBV	1683510	2421456	38.45%	2.966%
98	21.268	3103	3108	3125	rBV2	2433301	4423350	70.23%	5.419%
99	22.027	3234	3237	3247	rVB	1656976	2650344	42.08%	3.247%
100	23.986	3564	3570	3577	rBV	1434205	3341770	53.06%	4.094%

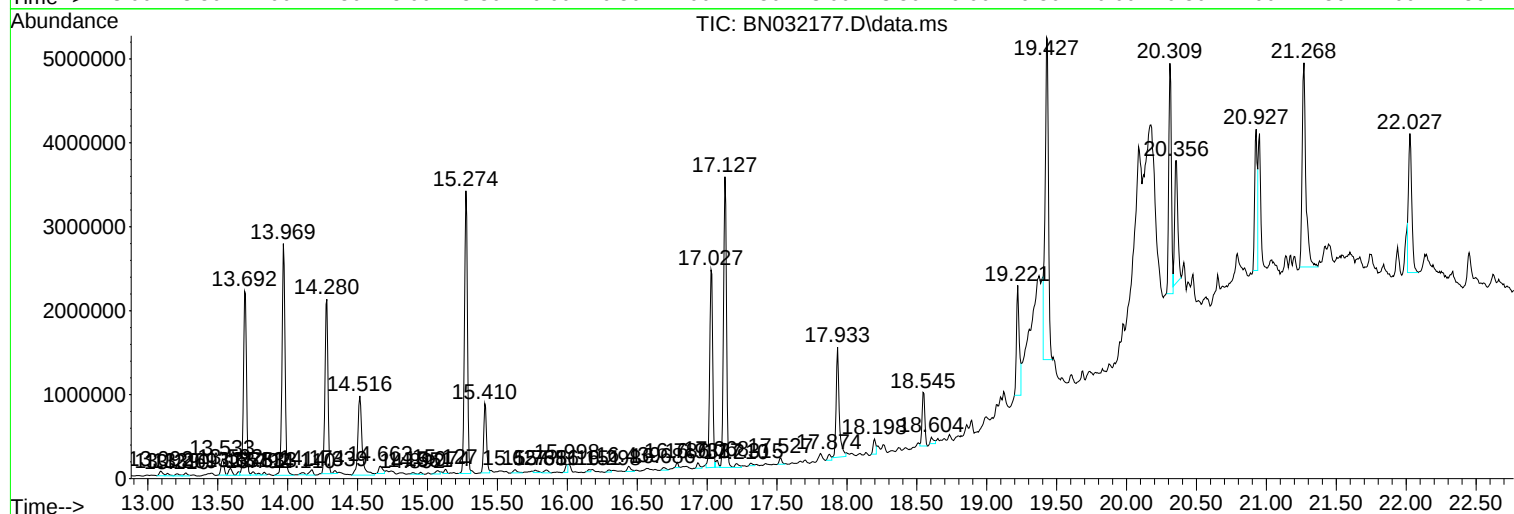
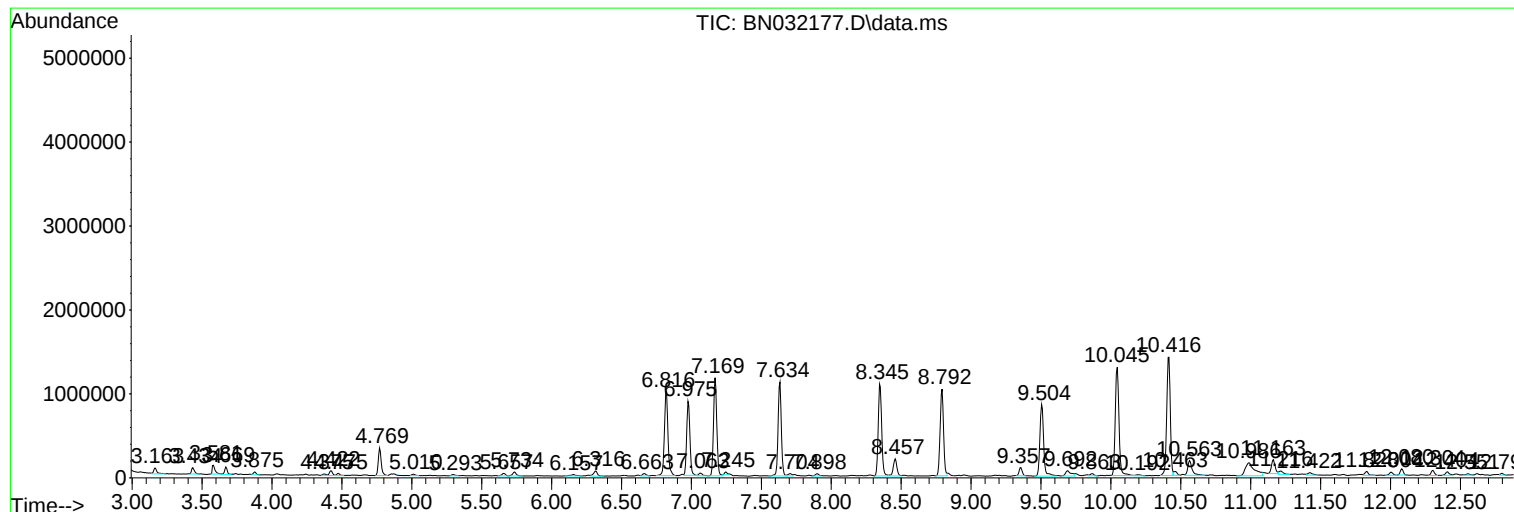
Sum of corrected areas: 81629107

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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



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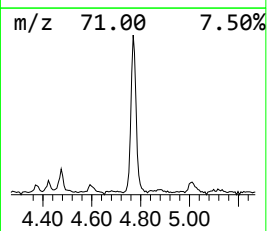
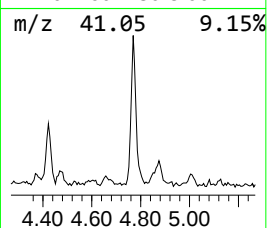
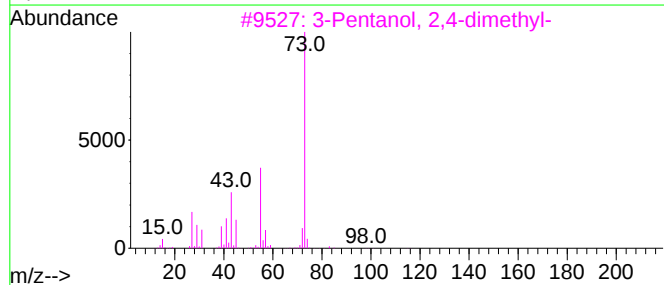
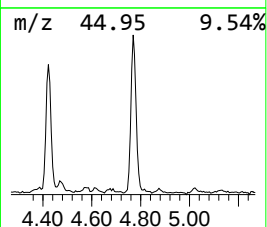
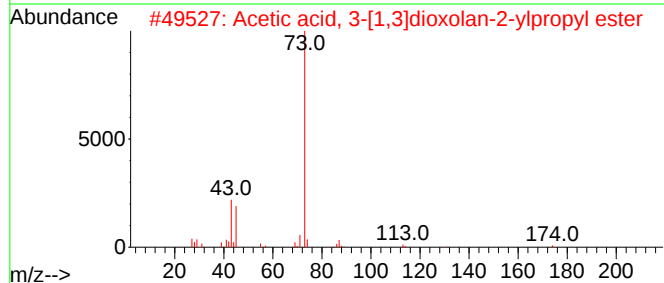
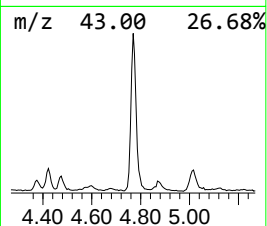
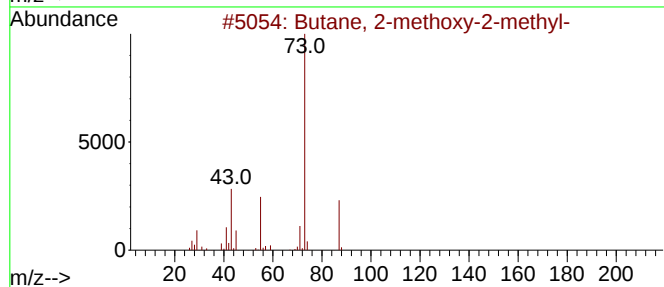
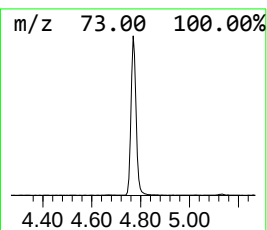
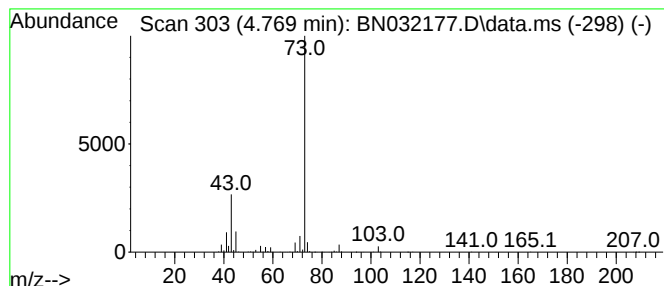
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN052924.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 unknown-01 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.769	5.31 ng/ul	486516	1,4-Dichlorobenzene-d4	7.634

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	38
2			Acetic acid, 3-[1,3]dioxolan-2-ylpropyl ester	174	C8H14O4	1000186-02-3	33
3			3-Pentanol, 2,4-dimethyl-	116	C7H16O	000600-36-2	33
4			Octanal, 7-methoxy-3,7-dimethyl-	186	C11H22O2	003613-30-7	28
5			o-2-Pentylhydroxylamine	103	C5H13NO	1000322-43-4	9



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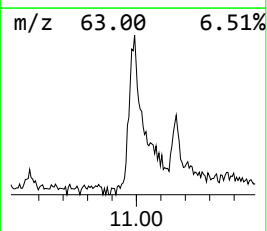
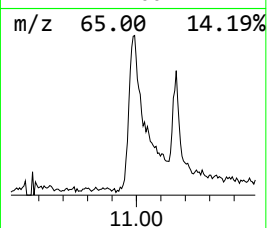
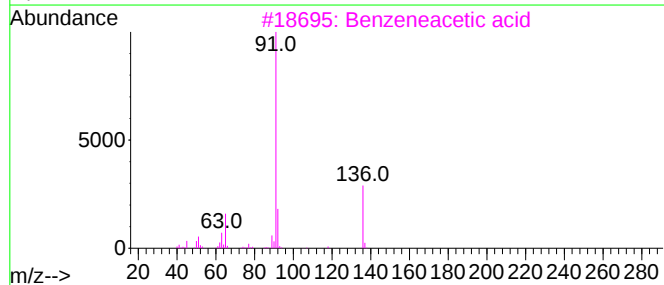
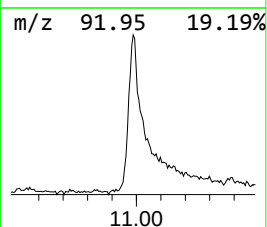
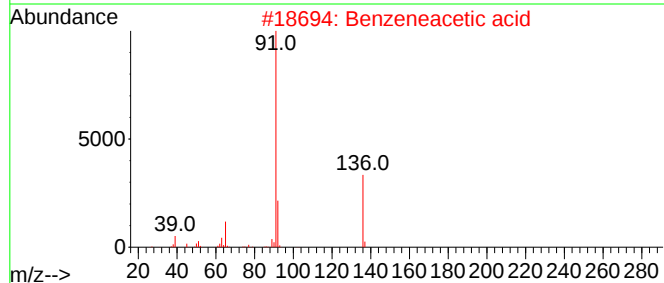
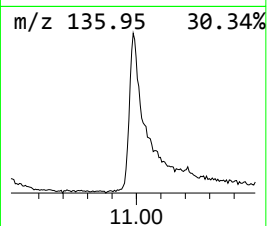
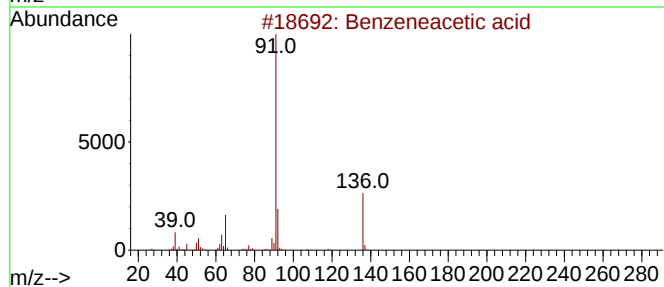
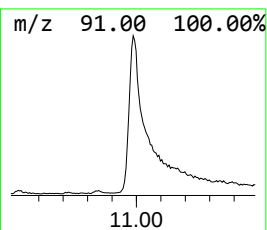
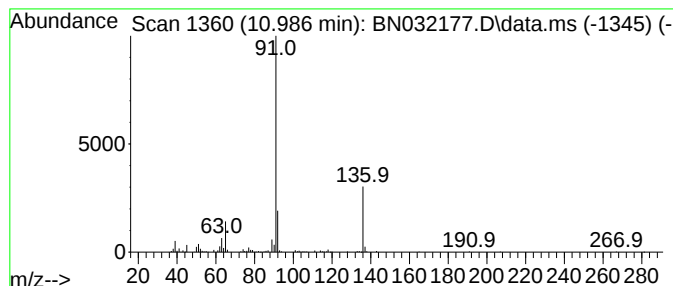
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN052924.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Benzeneacetic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.986	5.74 ng/ul	713629	Naphthalene-d8	10.416

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzeneacetic acid	136	C8H8O2	000103-82-2	91
2			Benzeneacetic acid	136	C8H8O2	000103-82-2	90
3			Benzeneacetic acid	136	C8H8O2	000103-82-2	81
4			Benzeneacetic acid	136	C8H8O2	000103-82-2	81
5			Benzeneacetic acid	136	C8H8O2	000103-82-2	52



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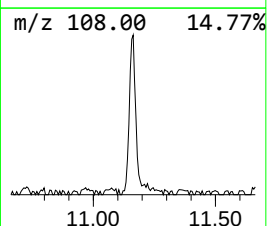
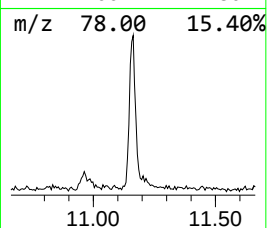
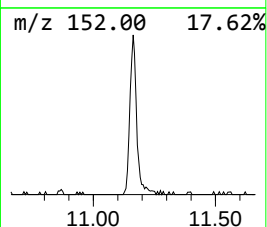
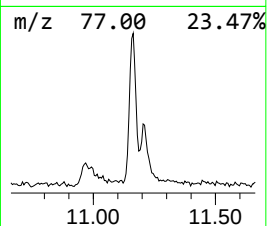
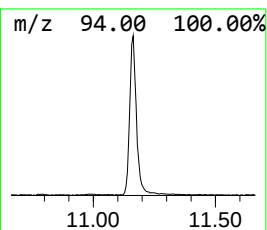
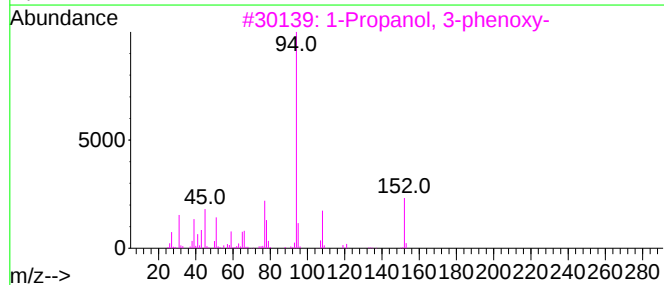
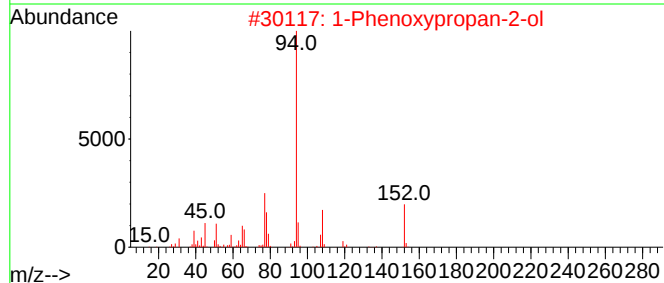
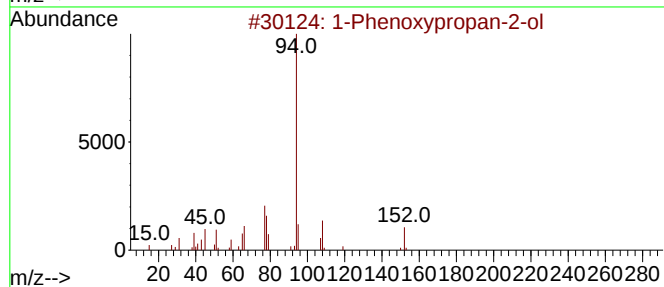
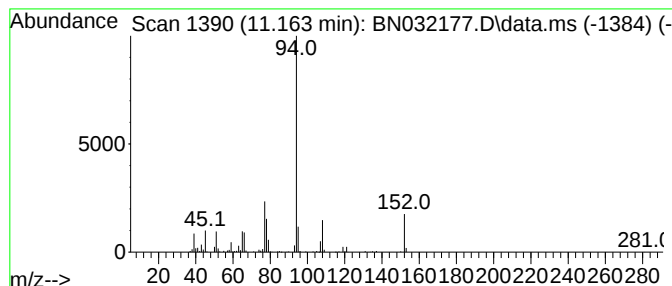
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN052924.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 1-Phenoxypropan-2-ol Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.163	2.38 ng/ul	295297	Naphthalene-d8	10.416

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	96
2			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	96
3			1-Propanol, 3-phenoxy-	152	C9H12O2	006180-61-6	91
4			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	90
5			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	70



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Data File : BN032177.D
Acq On : 22 Jun 2024 21:50
Operator : MA/JU
Sample : P2899-18
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
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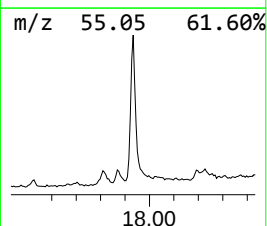
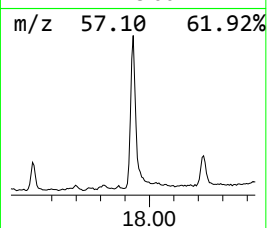
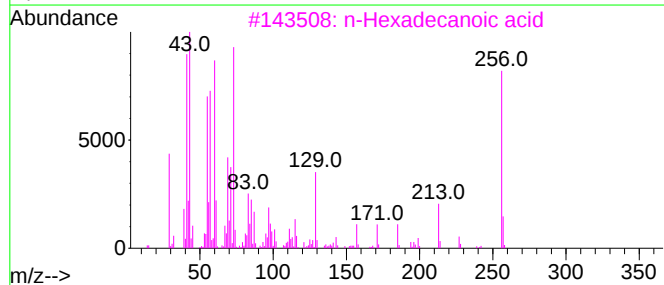
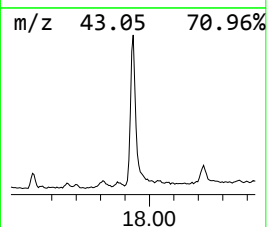
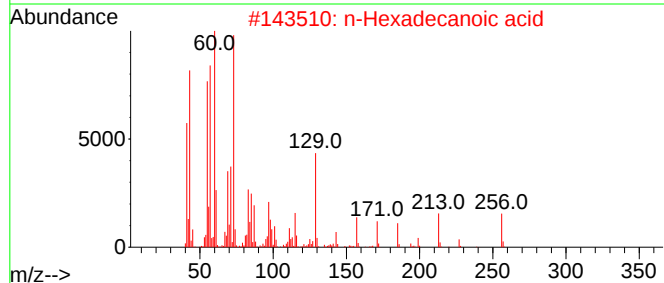
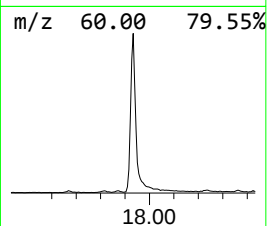
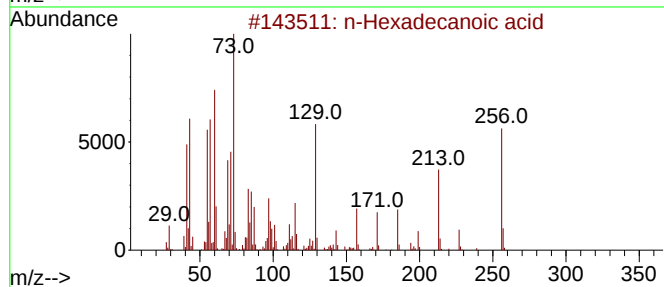
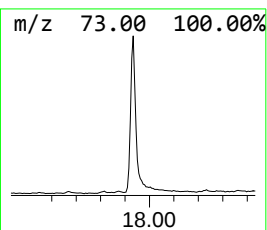
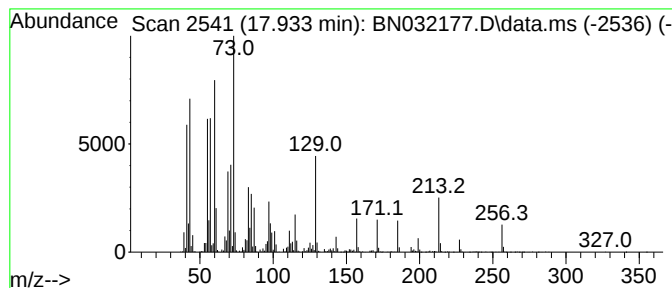
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.933	11.44 ng/ul	2009180	Phenanthrene-d10	17.027

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	96
5			Tridecanoic acid	214	C13H26O2	000638-53-9	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062224\
Data File : BN032177.D
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Sample : P2899-18
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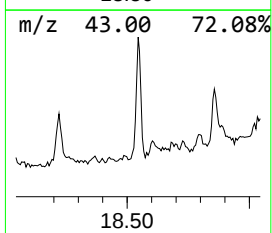
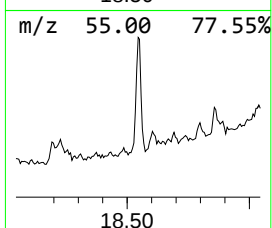
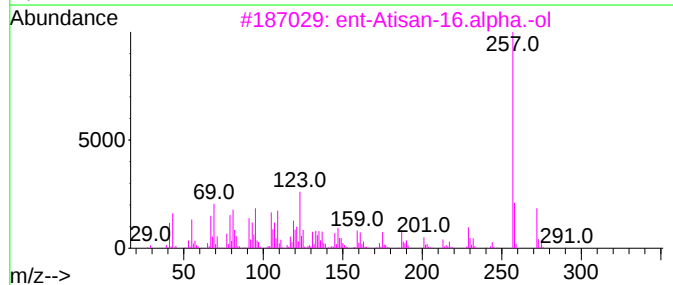
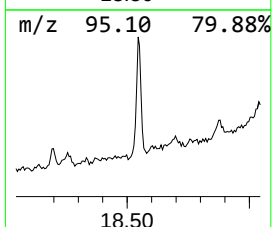
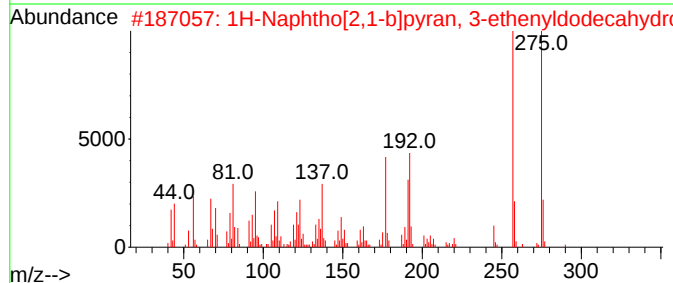
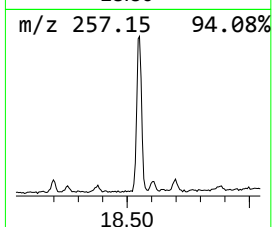
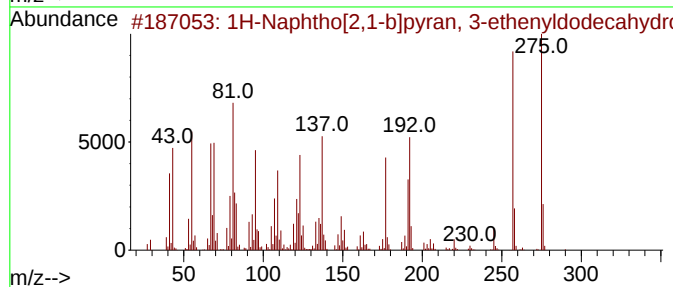
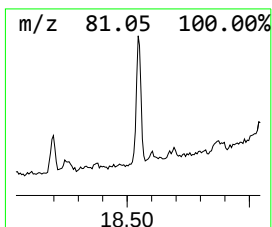
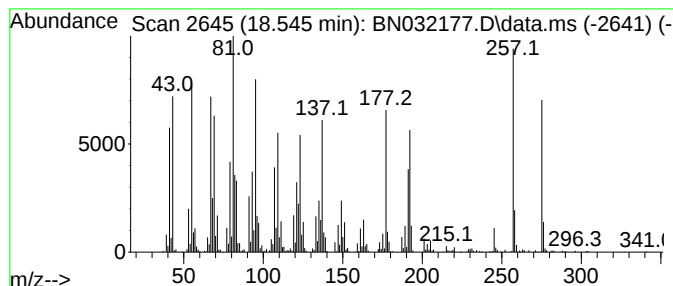
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 1H-Naphtho[2,1-b]pyran, 3-e... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.545	4.85 ng/ul	852265	Phenanthrene-d10	17.027

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Naphtho[2,1-b]pyran, 3-etheny...	290	C20H34O	000596-84-9	96
2			1H-Naphtho[2,1-b]pyran, 3-etheny...	290	C20H34O	000596-84-9	90
3			ent-Atisan-16.alpha.-ol	290	C20H34O	060801-64-1	55
4			(E)-3-Methyl-5-((1R,4aR,8aR)-5,5...	290	C20H34O	021738-29-4	55
5			1H-Naphtho[2,1-b]pyran, 3-etheny...	290	C20H34O	000596-84-9	49



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062224\
Data File : BN032177.D
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Sample : P2899-18
Misc :
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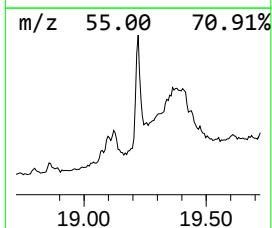
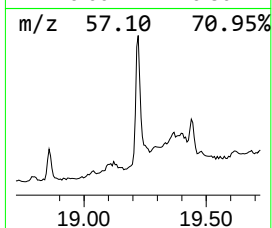
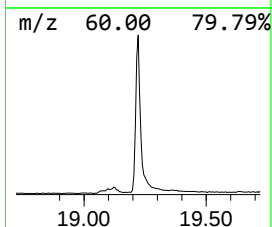
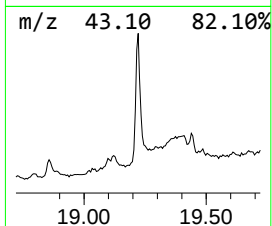
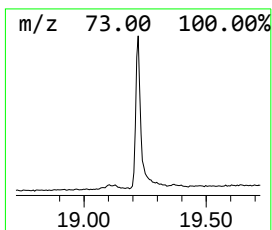
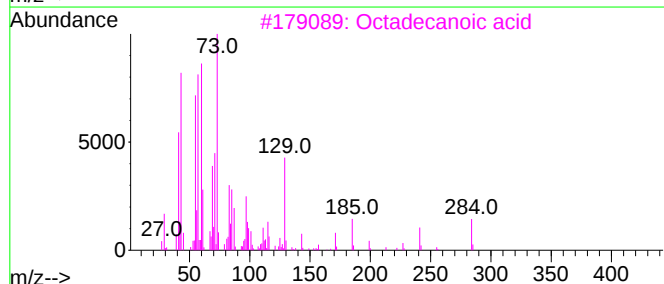
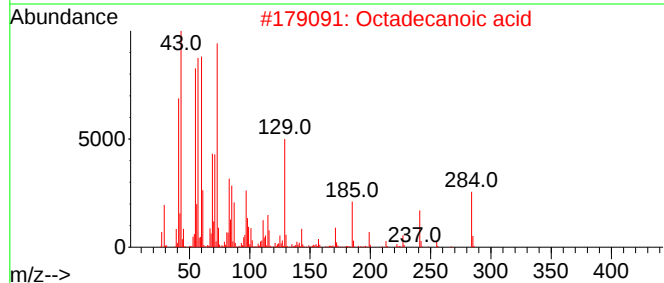
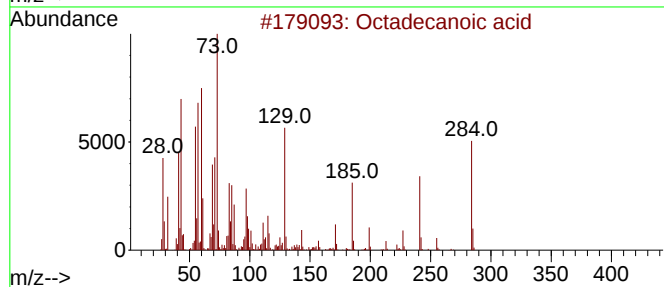
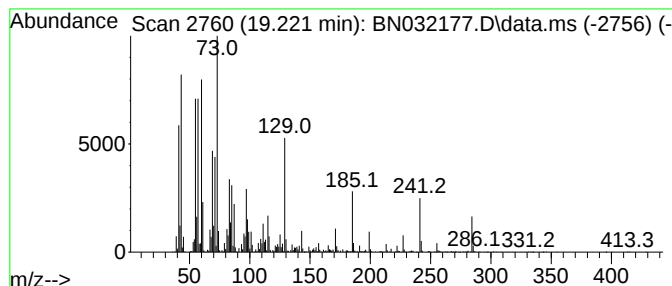
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Octadecanoic acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.221	8.62 ng/ul	1905700	Chrysene-d12	21.268

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	96
5			Octadecanoic acid	284	C18H36O2	000057-11-4	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062224\
Data File : BN032177.D
Acq On : 22 Jun 2024 21:50
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Sample : P2899-18
Misc :
ALS Vial : 19 Sample Multiplier: 1

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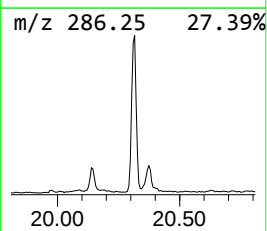
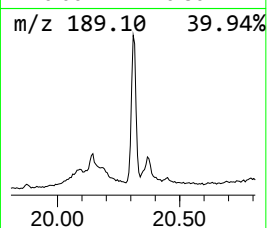
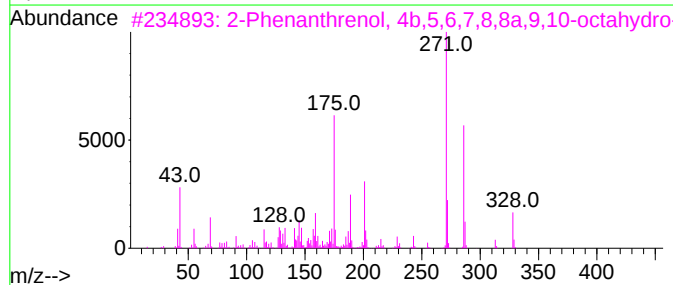
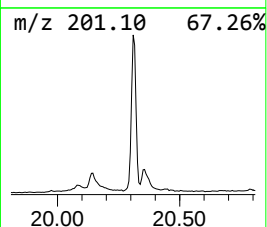
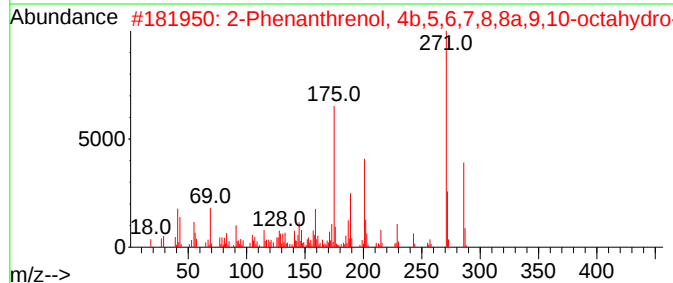
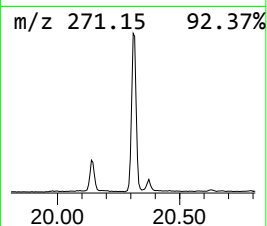
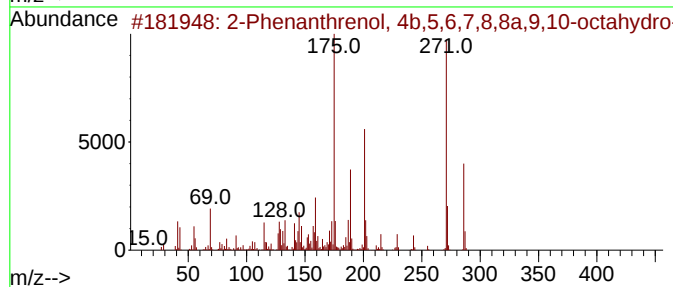
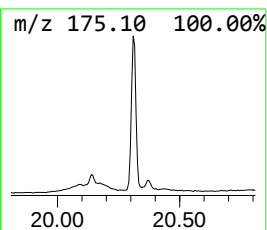
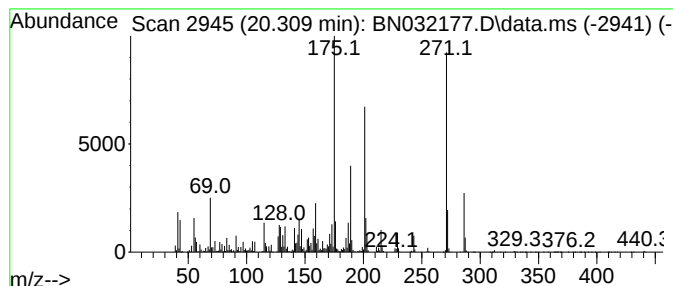
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 2-Phenanthrenol, 4b,5,6,7,8... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.309	16.22 ng/ul	3587000	Chrysene-d12	21.268

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Phenanthrenol, 4b,5,6,7,8,8a,9...	286	C20H30O	000511-15-9	99		
2	2-Phenanthrenol, 4b,5,6,7,8,8a,9...	286	C20H30O	000511-15-9	96		
3	2-Phenanthrenol, 4b,5,6,7,8,8a,9...	328	C22H32O2	015340-82-6	90		
4	2-Phenanthrenol, 4b,5,6,7,8,8a,9...	328	C22H32O2	015340-82-6	87		
5	13-Isopropylpodocarpene-12-ol-20-al	300	C20H28O2	024035-37-8	49		



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062224\
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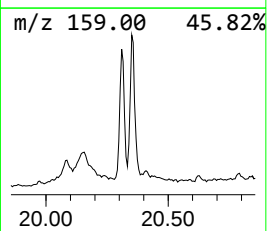
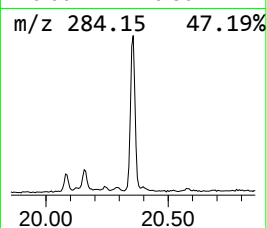
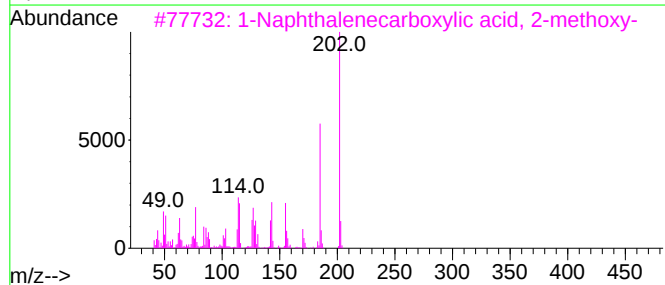
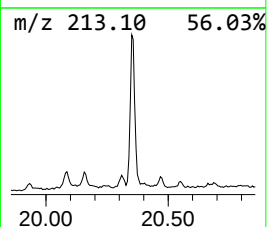
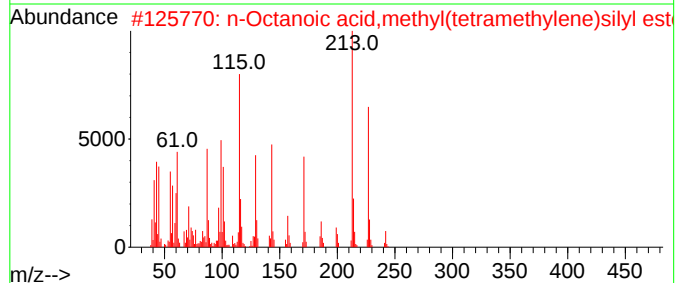
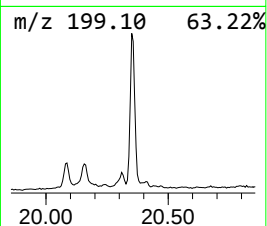
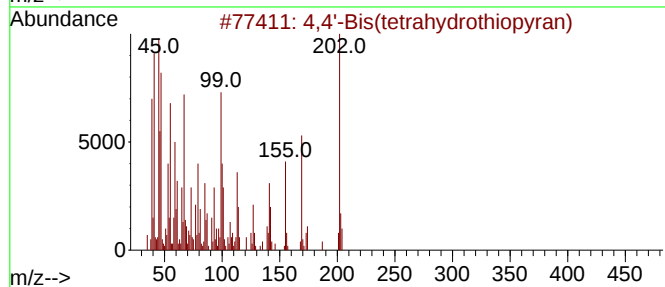
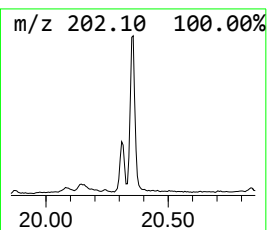
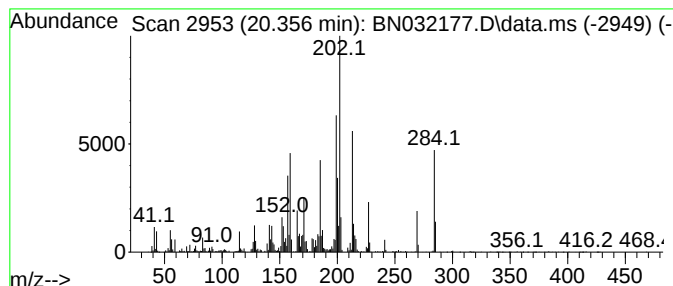
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 4,4'-Bis(tetrahydrothiopyran) Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.357	10.18 ng/ul	2252110	Chrysene-d12	21.268

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4,4'-Bis(tetrahydrothiopyran)	202	C10H18S2	116196-83-9	91
2			n-Octanoic acid,methyl(tetrameth...	242	C13H26O2Si	1010217-03-7	56
3			1-Naphthalenecarboxylic acid, 2-...	202	C12H10O3	000947-62-6	30
4			[1,2,5]-Oxadiazolo[3,4-a]cinnoli...	202	C10H10N4O	1000260-59-4	30
5			8-Amino-2,5-dimethyl-6-methoxyqu...	202	C12H14N2O	1000214-69-9	22



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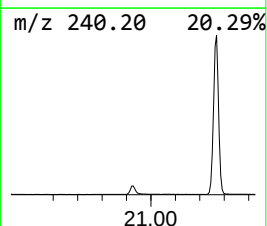
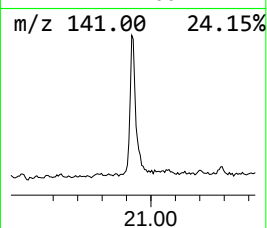
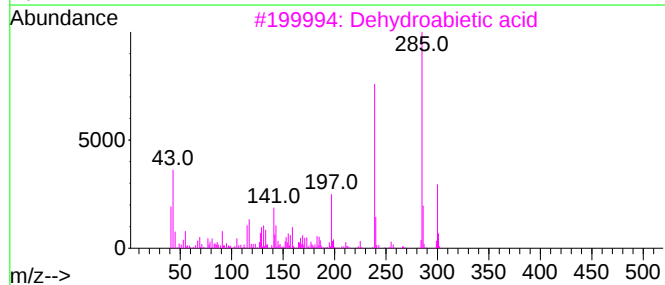
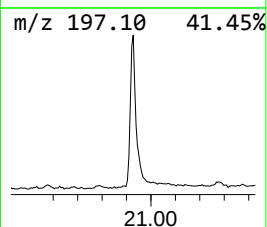
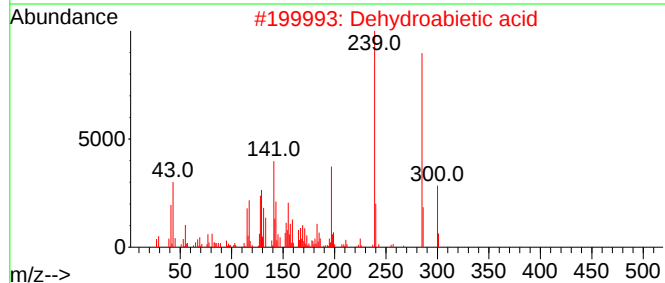
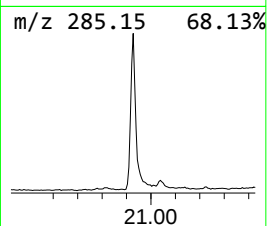
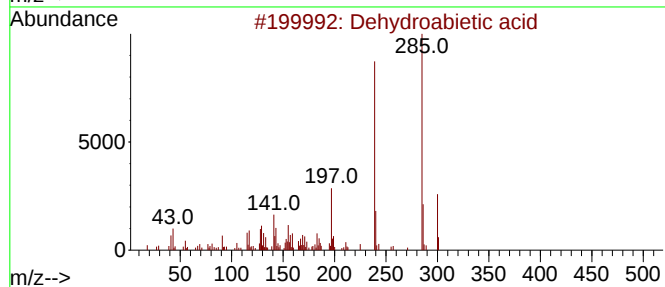
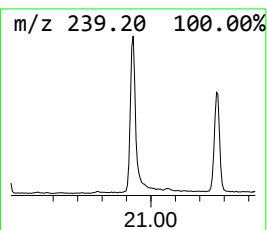
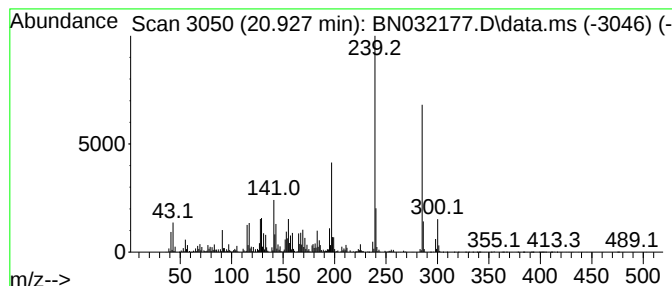
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 Dehydroabietic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.927	10.95 ng/ul	2421460	Chrysene-d12	21.268

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dehydroabietic acid	300	C20H28O2	001740-19-8	99
2			Dehydroabietic acid	300	C20H28O2	001740-19-8	99
3			Dehydroabietic acid	300	C20H28O2	001740-19-8	94
4			Dehydroabietic acid	300	C20H28O2	001740-19-8	91
5			trans-1,2-Bis(methyldichlorosilyl...	252	C4H8Cl4Si2	065899-10-7	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062224\
Data File : BN032177.D
Acq On : 22 Jun 2024 21:50
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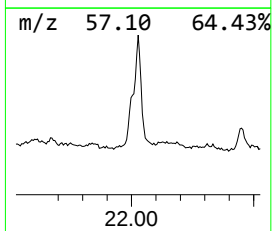
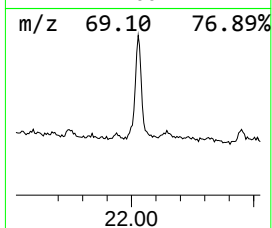
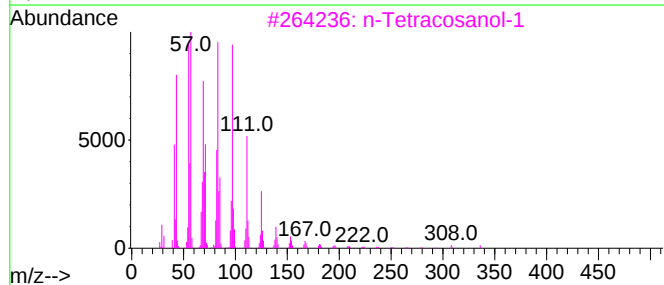
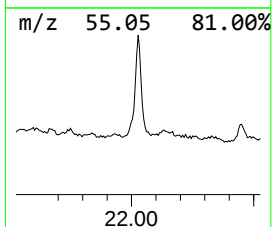
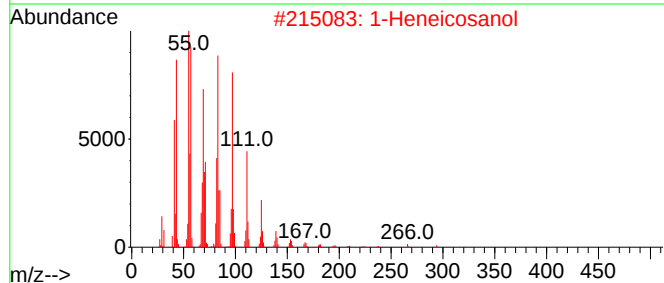
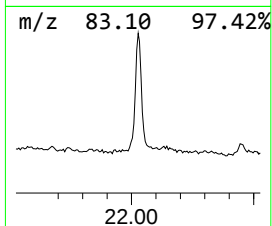
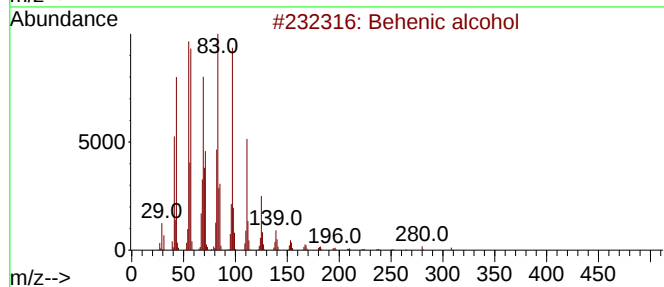
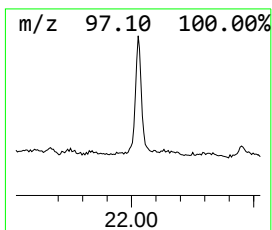
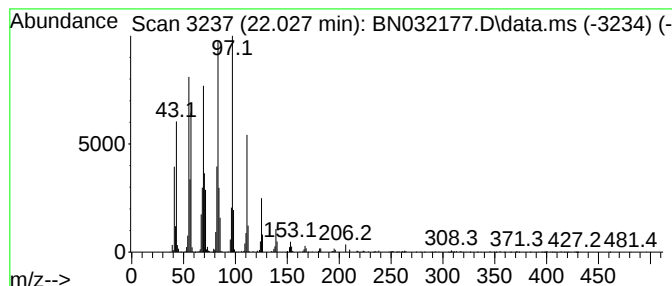
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN052924.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 10 Behenic alcohol Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.027	11.98 ng/ul	2650340	Chrysene-d12	21.268

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Behenic alcohol	326	C22H46O	000661-19-8	91
2			1-Heneicosanol	312	C21H44O	015594-90-8	91
3			n-Tetracosanol-1	354	C24H50O	000506-51-4	91
4			Cyclopentadecane	210	C15H30	000295-48-7	90
5			1-Hexacosene	364	C26H52	018835-33-1	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062224\
Data File : BN032177.D
Acq On : 22 Jun 2024 21:50
Operator : MA/JU
Sample : P2899-18
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
C0AB9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN052924.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
unknown-01	4.769	5.3	ng/ul	486516	1	7.634	1833070	20.0
Benzeneacetic acid	10.986	5.7	ng/ul	713629	2	10.416	2484510	20.0
1-Phenoxypropan...	11.163	2.4	ng/ul	295297	2	10.416	2484510	20.0
n-Hexadecanoic ...	17.933	11.4	ng/ul	2009180	4	17.027	3511890	20.0
1H-Naphtho[2,1-...	18.545	4.8	ng/ul	852265	4	17.027	3511890	20.0
Octadecanoic acid	19.221	8.6	ng/ul	1905700	5	21.268	4423350	20.0
2-Phenanthrenol...	20.309	16.2	ng/ul	3587000	5	21.268	4423350	20.0
4,4'-Bis(tetrah...	20.357	10.2	ng/ul	2252110	5	21.268	4423350	20.0
Dehydroabietic ...	20.927	10.9	ng/ul	2421460	5	21.268	4423350	20.0
Behenic alcohol	22.027	12.0	ng/ul	2650340	5	21.268	4423350	20.0