

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081524\
 Data File : BN033404.D
 Acq On : 15 Aug 2024 17:35
 Operator : MA/JU
 Sample : P3481-19
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 E0AF4

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-BN081424.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BN033404.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	40	rVB	123370	155827	2.17%	0.149%
2	3.416	68	73	85	rBV	429084	578891	8.06%	0.554%
3	3.552	91	96	110	rBV	1023132	1430953	19.93%	1.370%
4	3.869	145	150	156	rVB	30866	42293	0.59%	0.040%
5	4.510	256	259	268	rVB6	10647	20008	0.28%	0.019%
6	4.975	332	338	343	rBV3	13672	23193	0.32%	0.022%
7	5.699	455	461	467	rBV	57327	83279	1.16%	0.080%
8	6.516	596	600	605	rBV3	17994	26109	0.36%	0.025%
9	6.581	605	611	619	rVB2	19054	34005	0.47%	0.033%
10	6.751	632	640	654	rVB	1822309	2821850	39.30%	2.701%
11	6.875	656	661	662	rBV2	17189	23651	0.33%	0.023%
12	6.916	662	668	676	rVB	1529748	2258924	31.46%	2.162%
13	7.104	692	700	709	rBV	1836895	2863525	39.88%	2.741%
14	7.181	709	713	724	rVB	191804	284388	3.96%	0.272%
15	7.569	771	779	787	rVV	2106543	3205990	44.65%	3.069%
16	7.646	787	792	798	rVB	44628	63051	0.88%	0.060%
17	7.869	827	830	836	rVV5	8265	14924	0.21%	0.014%
18	7.934	838	841	846	rVB4	8910	14184	0.20%	0.014%
19	8.275	891	899	907	rBV	2046658	3254838	45.33%	3.115%
20	8.710	959	973	984	rBV	1588485	2541241	35.39%	2.432%
21	8.840	991	995	999	rBV4	9024	13884	0.19%	0.013%
22	8.893	1000	1004	1008	rVB5	9052	13349	0.19%	0.013%
23	9.057	1025	1032	1036	rBV2	18184	29233	0.41%	0.028%
24	9.193	1050	1055	1060	rVB2	21145	35432	0.49%	0.034%
25	9.287	1064	1071	1078	rBV	171791	257110	3.58%	0.246%
26	9.422	1086	1094	1104	rBV	1172093	1888982	26.31%	1.808%
27	9.669	1132	1136	1141	rVB5	9244	14388	0.20%	0.014%
28	9.745	1144	1149	1155	rBV4	17608	34690	0.48%	0.033%
29	9.898	1170	1175	1178	rVV	13804	22441	0.31%	0.021%
30	9.957	1178	1185	1197	rVB	1944686	3110700	43.32%	2.977%
31	10.140	1209	1216	1222	rVB8	10434	18009	0.25%	0.017%
32	10.334	1240	1249	1257	rBV	2525134	4163030	57.98%	3.985%
33	10.469	1264	1272	1286	rBV	800310	1313867	18.30%	1.258%
34	10.892	1339	1344	1350	rVB	70782	102354	1.43%	0.098%
35	11.081	1367	1376	1382	rBV	375412	617085	8.59%	0.591%
36	11.134	1382	1385	1392	rVB8	11910	19835	0.28%	0.019%

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Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN081424.MA.M
Title : SVOA CALIBRATION

37	11.251	1397	1405	1409	rBV9	7570	17710	0.25%	0.017%
38	11.928	1516	1520	1526	rVB2	40293	63853	0.89%	0.061%
39	12.822	1669	1672	1676	rBV4	9524	13837	0.19%	0.013%
40	13.063	1705	1713	1718	rBV3	17387	26452	0.37%	0.025%

41	13.539	1785	1794	1799	rBV6	25717	59219	0.82%	0.057%
42	13.628	1803	1809	1815	rVV	2748969	3751516	52.25%	3.591%
43	13.681	1815	1818	1821	rVV5	14069	19351	0.27%	0.019%
44	13.763	1829	1832	1836	rVV4	9832	13033	0.18%	0.012%
45	13.898	1847	1855	1865	rBV	3108964	4858467	67.67%	4.650%

46	14.069	1879	1884	1891	rVB9	9853	20006	0.28%	0.019%
47	14.210	1901	1908	1916	rVV	3149149	4676773	65.14%	4.476%
48	14.269	1916	1918	1926	rVB6	15591	24809	0.35%	0.024%
49	14.410	1932	1942	1945	rBV	1151577	1670004	23.26%	1.598%
50	14.445	1945	1948	1949	rVV	514873	655342	9.13%	0.627%

51	14.463	1949	1951	1962	rVB	607862	682427	9.50%	0.653%
52	14.733	1993	1997	2003	rBV4	9313	16008	0.22%	0.015%
53	15.045	2046	2050	2054	rBV4	17327	25464	0.35%	0.024%
54	15.110	2057	2061	2064	rBV	10409	13161	0.18%	0.013%
55	15.204	2071	2077	2084	rBV	3851056	5491738	76.49%	5.256%

56	15.257	2084	2086	2090	rVB3	13362	18487	0.26%	0.018%
57	15.322	2091	2097	2111	rVV	805961	1121170	15.62%	1.073%
58	15.445	2115	2118	2123	rVV2	22027	29603	0.41%	0.028%
59	15.786	2171	2176	2184	rVB5	28676	45100	0.63%	0.043%
60	15.975	2203	2208	2210	rBV3	16651	24043	0.33%	0.023%

61	15.998	2210	2212	2220	rVB7	15394	25176	0.35%	0.024%
62	16.392	2275	2279	2284	rVB5	18361	25840	0.36%	0.025%
63	16.551	2301	2306	2309	rBV7	11512	16947	0.24%	0.016%
64	16.627	2314	2319	2327	rVB	97663	140362	1.95%	0.134%
65	16.710	2329	2333	2336	rBV5	10660	17616	0.25%	0.017%

66	16.745	2336	2339	2341	rBV2	14689	16290	0.23%	0.016%
67	16.957	2367	2375	2380	rBV	3512639	4822083	67.16%	4.615%
68	16.992	2380	2381	2385	rVV	36894	42627	0.59%	0.041%
69	17.057	2385	2392	2402	rVB	3699510	5329342	74.23%	5.101%
70	17.163	2406	2410	2414	rVB6	27029	38901	0.54%	0.037%

71	17.345	2432	2441	2466	rBV	1892036	3970606	55.30%	3.800%
72	17.586	2478	2482	2488	rVB2	57733	83067	1.16%	0.080%
73	17.674	2494	2497	2502	rVB2	49281	64142	0.89%	0.061%
74	17.757	2508	2511	2519	rVB6	27803	46911	0.65%	0.045%
75	17.892	2528	2534	2542	rBV	443063	661944	9.22%	0.634%

76	18.180	2579	2583	2590	rVB4	43429	77807	1.08%	0.074%
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77	18.333	2606	2609	2615	rVB7	16657	23737	0.33%	0.023%
78	18.763	2677	2682	2688	rBV6	35053	66597	0.93%	0.064%
79	18.857	2694	2698	2702	rVB5	53789	69690	0.97%	0.067%
80	18.921	2705	2709	2715	rVB2	56857	77617	1.08%	0.074%
81	18.992	2716	2721	2724	rBV	70779	113384	1.58%	0.109%
82	19.021	2724	2726	2729	rVB	47587	49909	0.70%	0.048%
83	19.180	2749	2753	2758	rBV2	141231	192417	2.68%	0.184%
84	19.357	2777	2783	2789	rBV	4573565	6207145	86.45%	5.941%
85	19.421	2789	2794	2815	rVB	1697428	2780161	38.72%	2.661%
86	19.921	2876	2879	2882	rBV4	22037	27578	0.38%	0.026%
87	20.280	2937	2940	2944	rVB2	70499	81881	1.14%	0.078%
88	20.504	2973	2978	2983	rVB	603941	727708	10.14%	0.697%
89	20.674	3003	3007	3013	rBV6	116502	168268	2.34%	0.161%
90	20.792	3022	3027	3032	rBV2	1178207	1749535	24.37%	1.675%
91	20.851	3032	3037	3041	rVB4	238104	404658	5.64%	0.387%
92	20.921	3045	3049	3062	rVV2	633053	1189597	16.57%	1.139%
93	21.045	3066	3070	3077	rVB3	327958	524409	7.30%	0.502%
94	21.192	3087	3095	3104	rVB	3628934	5775001	80.43%	5.527%
95	21.986	3226	3230	3237	rVB3	199734	344397	4.80%	0.330%
96	22.762	3359	3362	3367	rVB	109522	145846	2.03%	0.140%
97	23.309	3450	3455	3460	rBV	207392	420627	5.86%	0.403%
98	23.833	3534	3544	3562	rVB	2928282	7179924	100.00%	6.872%
99	24.021	3567	3576	3587	rVB	2446973	5782019	80.53%	5.534%
100	25.127	3761	3764	3772	rVB2	121029	264447	3.68%	0.253%

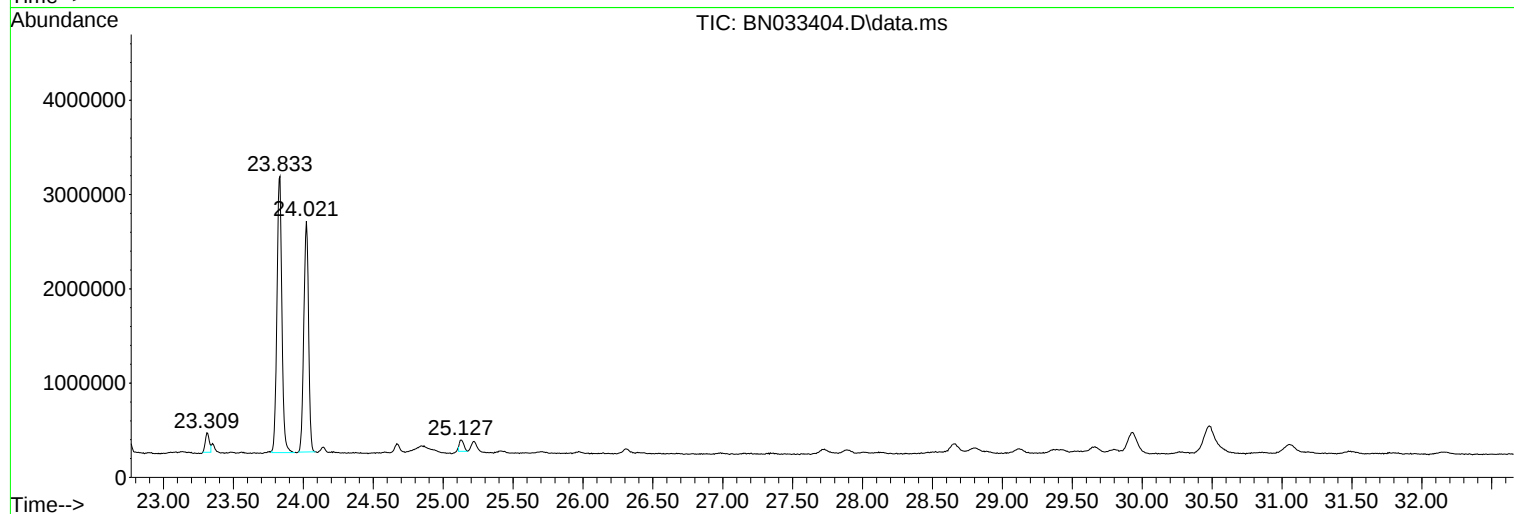
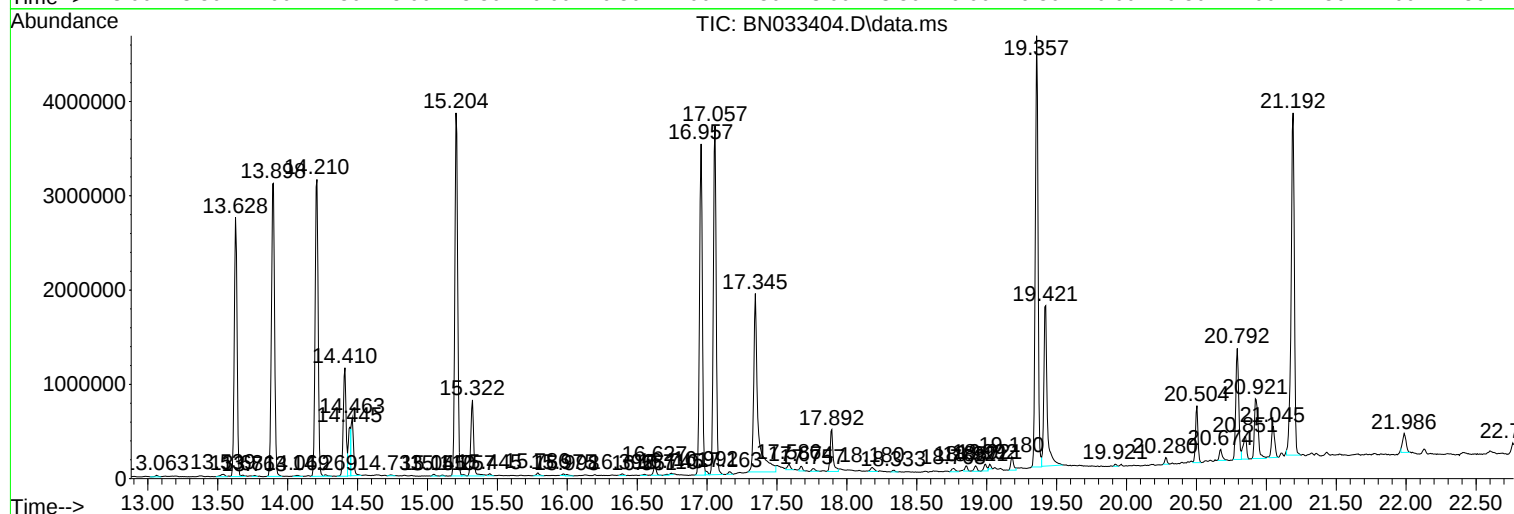
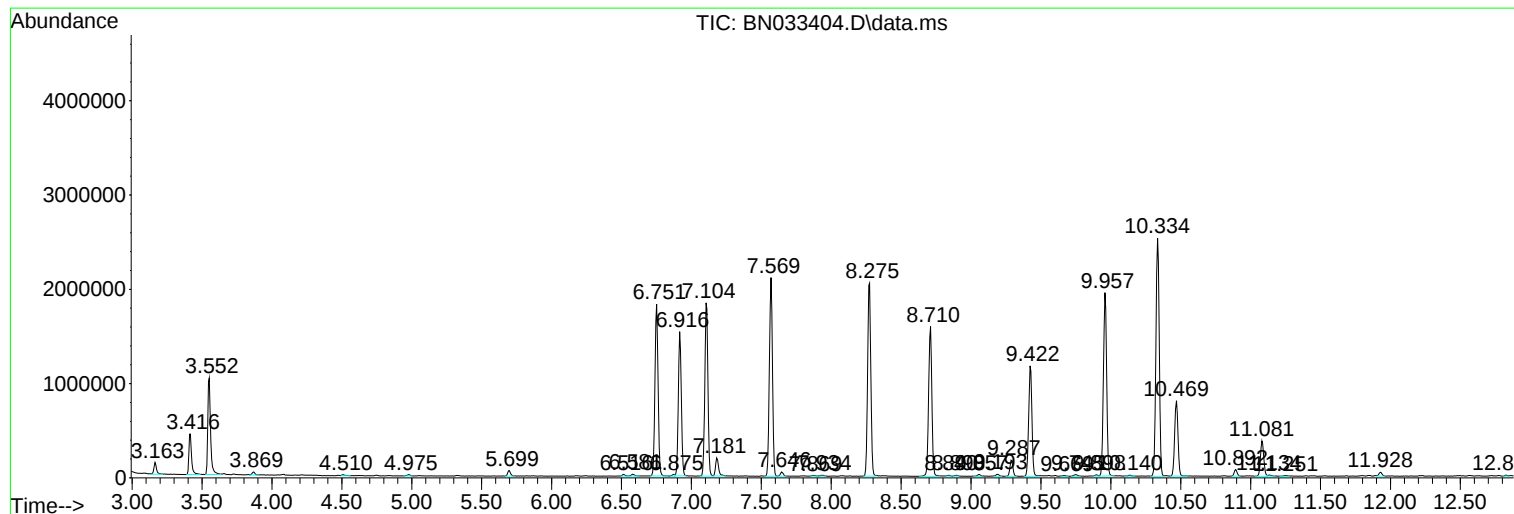
Sum of corrected areas: 104479299

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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN081424.MA.M
Quant Title : SVOA CALIBRATION

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



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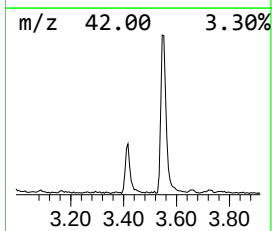
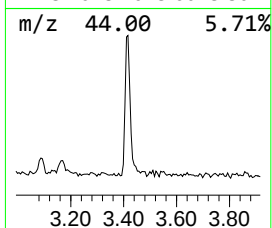
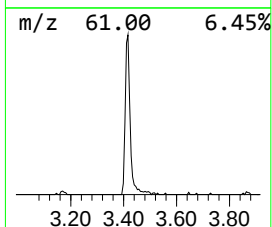
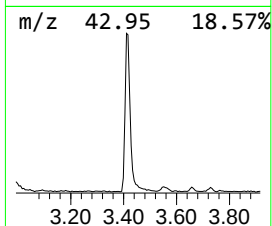
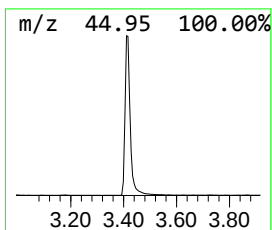
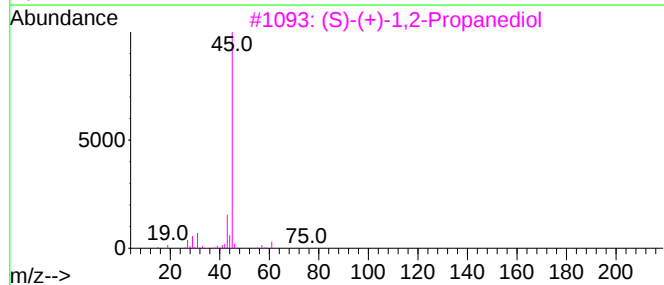
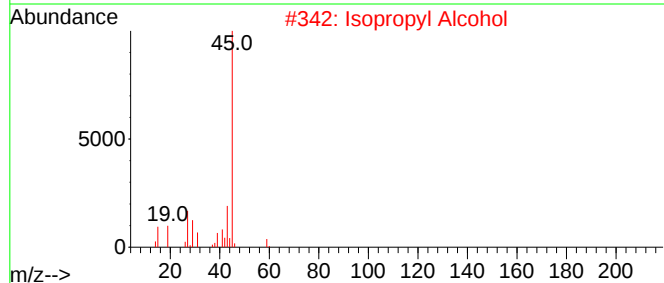
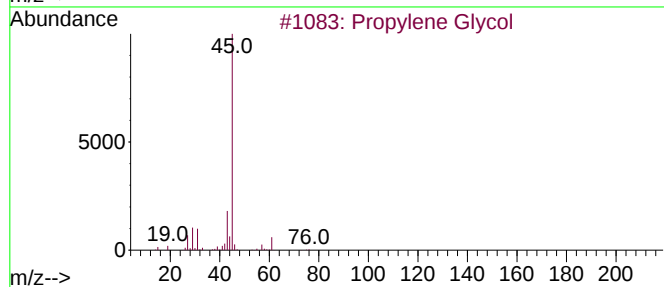
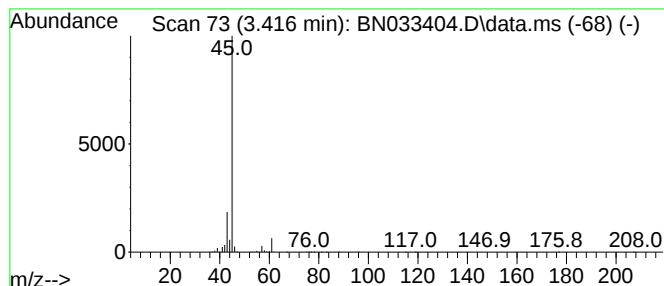
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN081424.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Propylene Glycol Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.416	3.61 ng/ul	578891	1,4-Dichlorobenzene-d4	7.569

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Propylene Glycol	76	C3H8O2	000057-55-6	90
2			Isopropyl Alcohol	60	C3H8O	000067-63-0	56
3			(S)-(+)-1,2-Propanediol	76	C3H8O2	004254-15-3	43
4			Propanoic acid, 2-hydroxy-, meth...	104	C4H8O3	000547-64-8	40
5			Isopropyl Alcohol	60	C3H8O	000067-63-0	9



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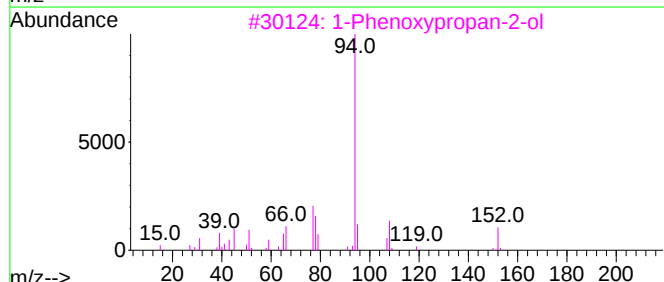
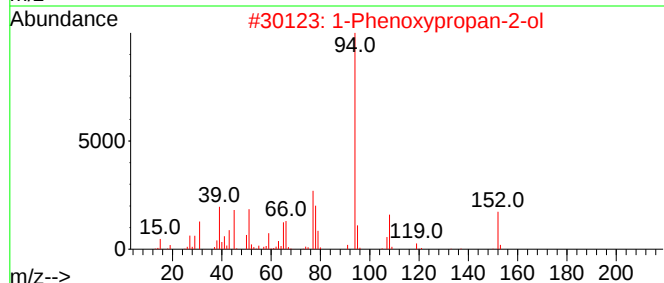
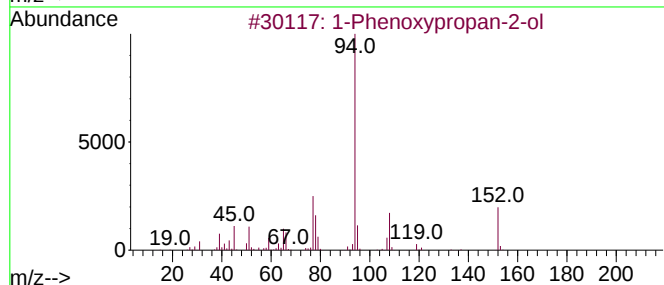
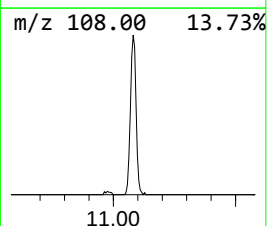
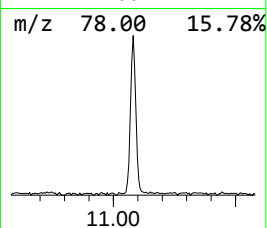
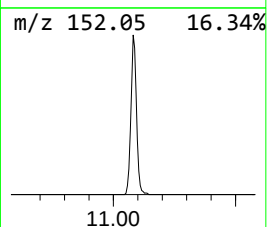
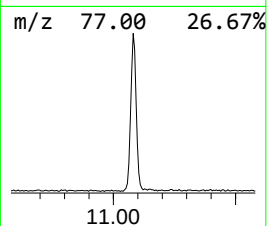
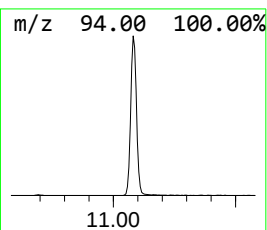
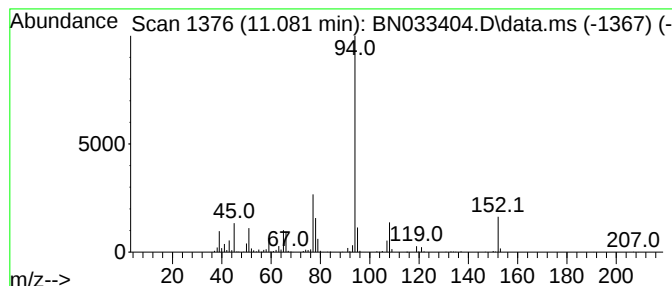
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN081424.MA.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.081	2.96 ng/ul	617085	Naphthalene-d8	10.334

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	95
3			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	94
4			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	94
5			1-Propanol, 3-phenoxy-	152	C9H12O2	006180-61-6	87



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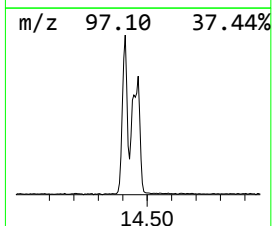
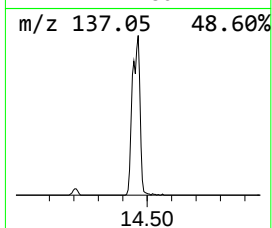
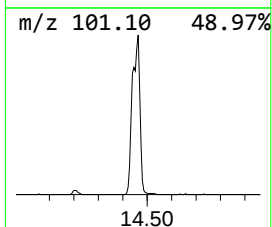
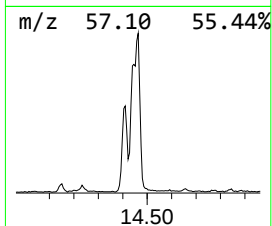
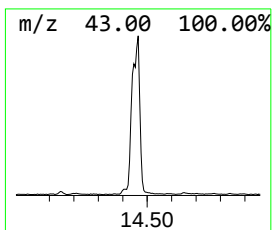
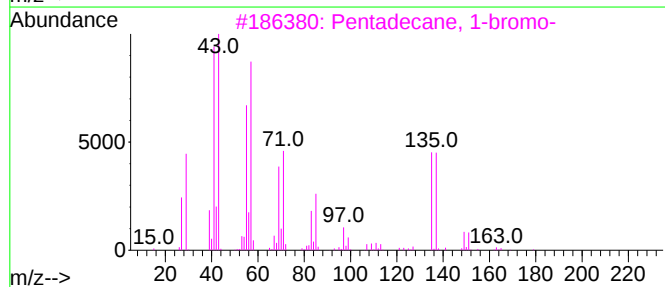
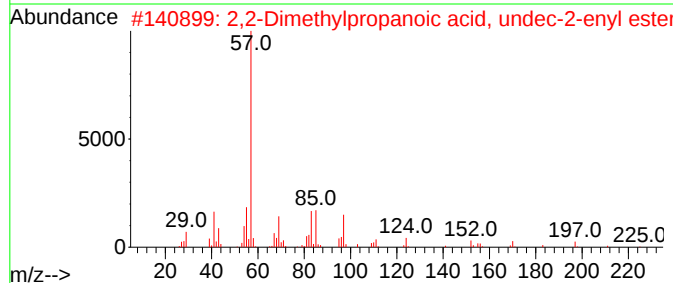
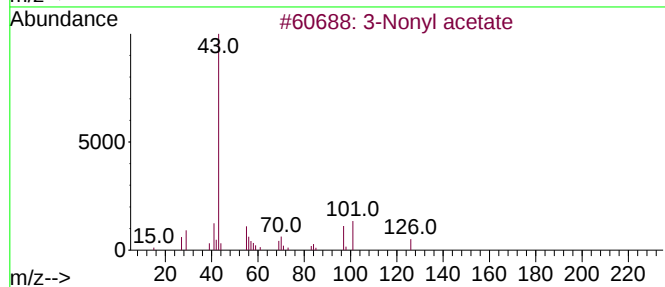
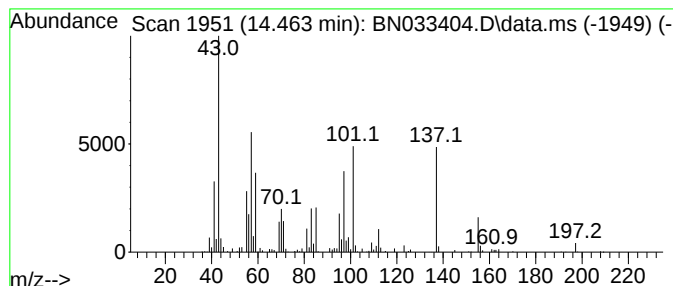
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 unknown-01 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.463	2.92 ng/ul	682427	Acenaphthene-d10	14.210

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Nonyl acetate	186	C11H22O2	1010429-16-2	25
2			2,2-Dimethylpropanoic acid, unde...	254	C16H30O2	1000299-33-7	14
3			Pentadecane, 1-bromo-	290	C15H31Br	000629-72-1	12
4			Nonane, 1-bromo-	206	C9H19Br	000693-58-3	12
5			Carbonic acid, prop-1-en-2-yl un...	256	C15H28O3	1000382-90-6	10



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081524\
Data File : BN033404.D
Acq On : 15 Aug 2024 17:35
Operator : MA/JU
Sample : P3481-19
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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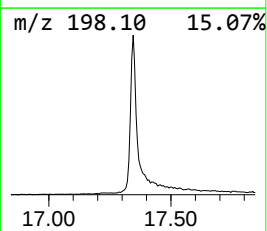
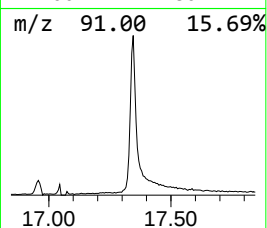
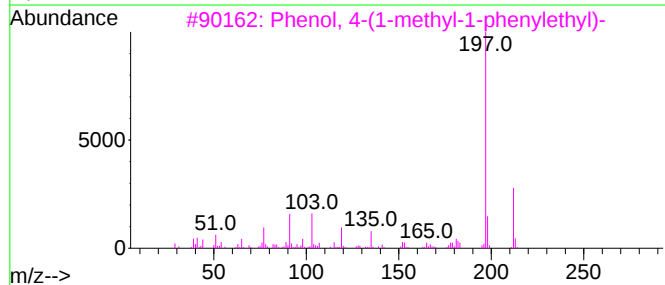
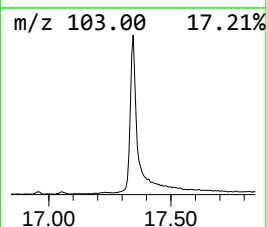
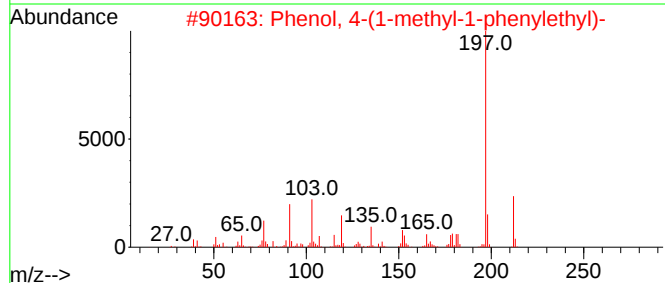
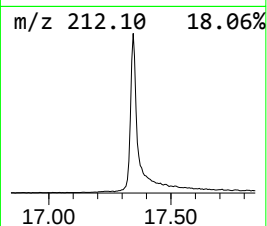
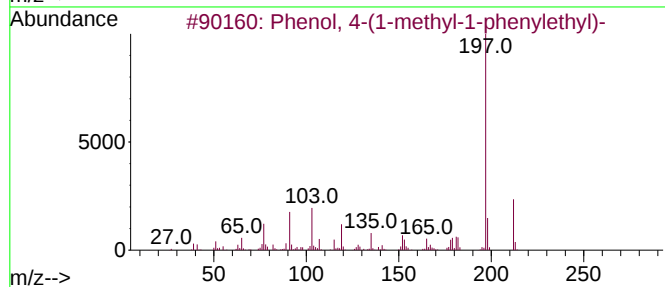
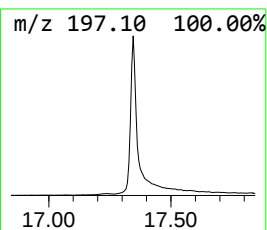
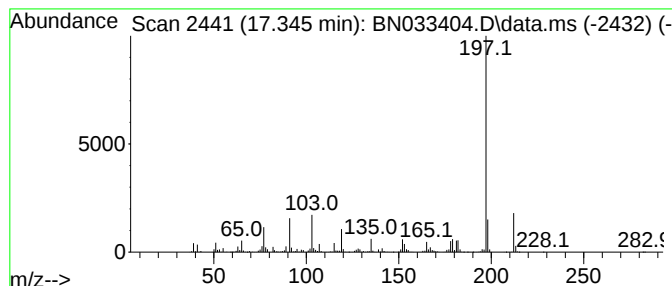
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Phenol, 4-(1-methyl-1-pheny... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.345	16.47 ng/ul	3970610	Phenanthrene-d10	16.957

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenol, 4-(1-methyl-1-phenylethyl)-	212	C15H16O	000599-64-4	98
2			Phenol, 4-(1-methyl-1-phenylethyl)-	212	C15H16O	000599-64-4	97
3			Phenol, 4-(1-methyl-1-phenylethyl)-	212	C15H16O	000599-64-4	91
4			2,3-Dimethyl-4-biphenylamine	197	C14H15N	022750-87-4	70
5			Dimethyl diphenylsilane	212	C14H16Si	000778-24-5	68



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081524\
Data File : BN033404.D
Acq On : 15 Aug 2024 17:35
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Sample : P3481-19
Misc :
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Instrument :
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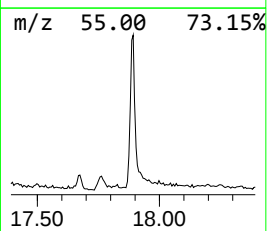
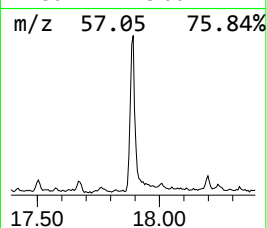
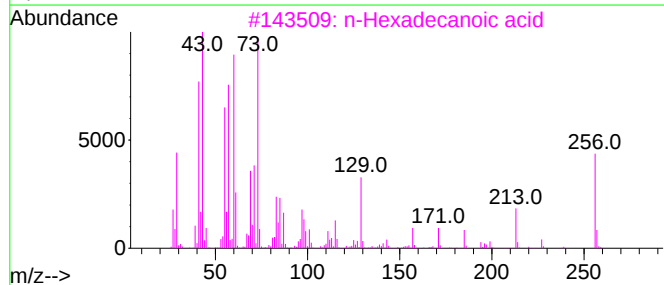
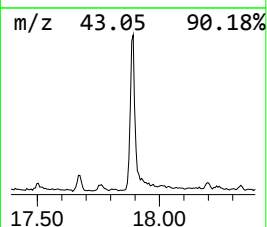
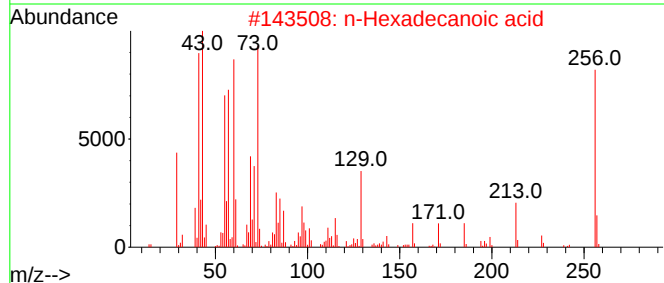
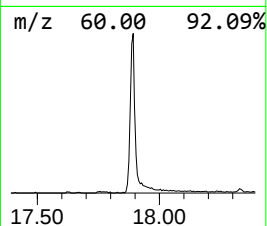
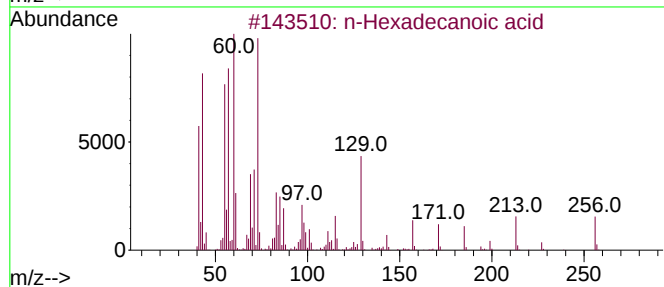
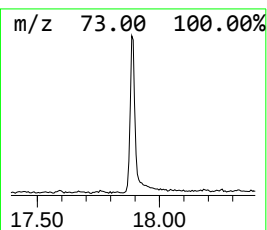
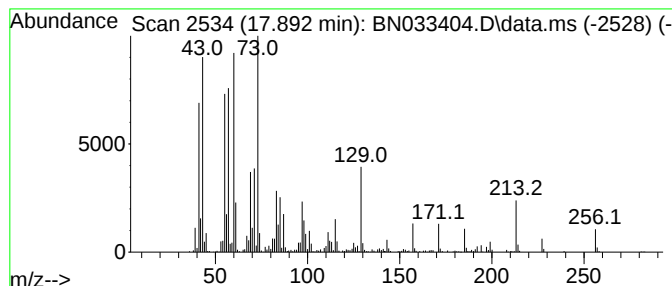
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 n-Hexadecanoic acid Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.892	2.75 ng/ul	661944	Phenanthrene-d10	16.957

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
3			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
5			Tridecanoic acid	214	C13H26O2	000638-53-9	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081524\
Data File : BN033404.D
Acq On : 15 Aug 2024 17:35
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Sample : P3481-19
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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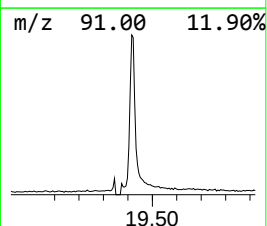
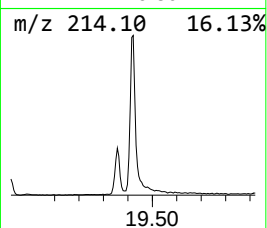
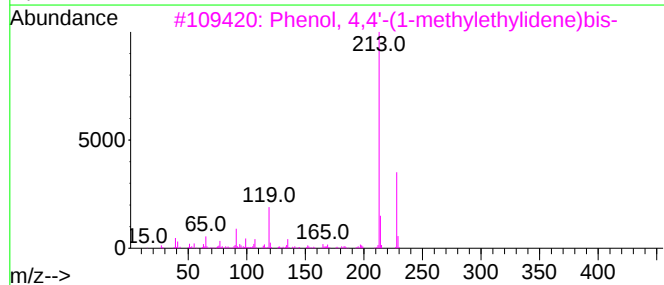
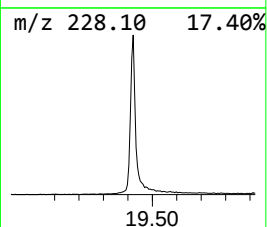
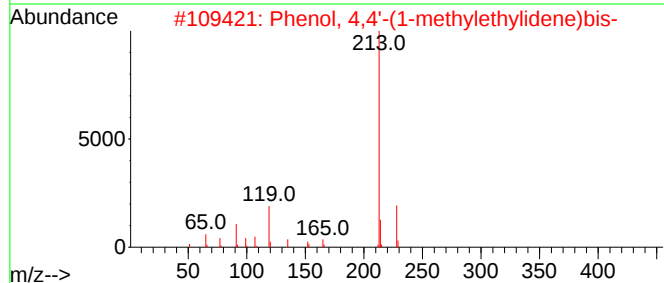
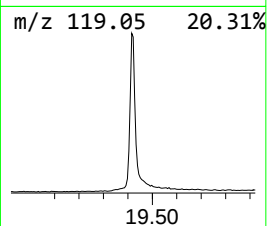
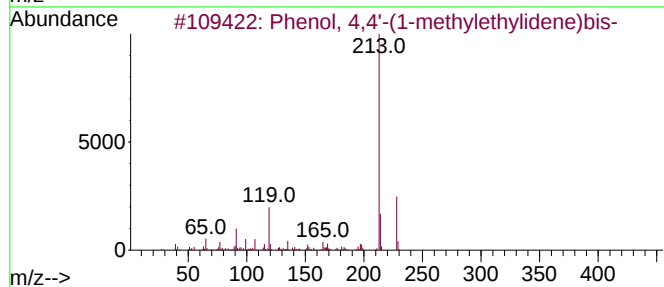
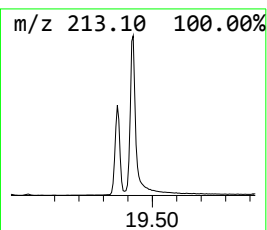
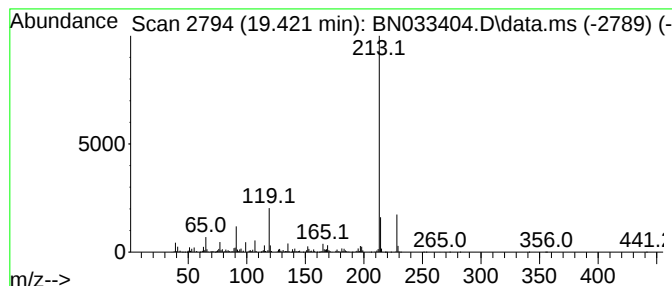
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Phenol, 4,4'-(1-methylethyl... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.421	9.63 ng/ul	2780160	Chrysene-d12	21.192

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenol, 4,4'-(1-methylethylidene...	228	C15H16O2	000080-05-7	98
2			Phenol, 4,4'-(1-methylethylidene...	228	C15H16O2	000080-05-7	98
3			Phenol, 4,4'-(1-methylethylidene...	228	C15H16O2	000080-05-7	91
4			Phenol, 4,4'-(1-methylethylidene...	228	C15H16O2	000080-05-7	91
5			4,4'-(1,3-Dimethylbutylidene)bis...	270	C18H22O2	006807-17-6	80



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081524\
Data File : BN033404.D
Acq On : 15 Aug 2024 17:35
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Sample : P3481-19
Misc :
ALS Vial : 13 Sample Multiplier: 1

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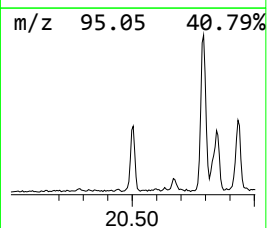
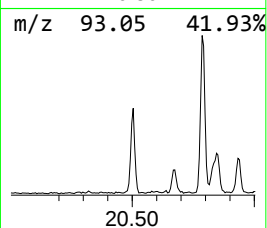
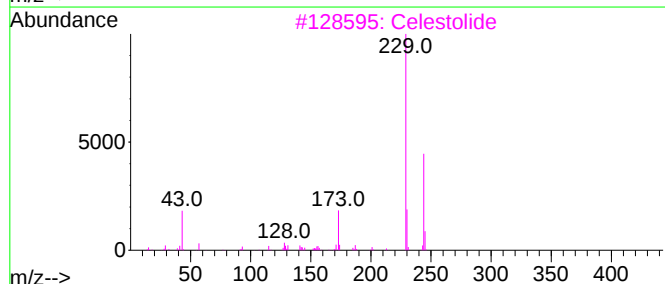
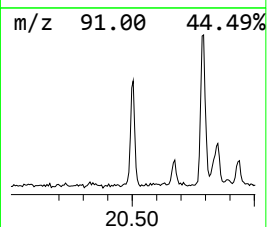
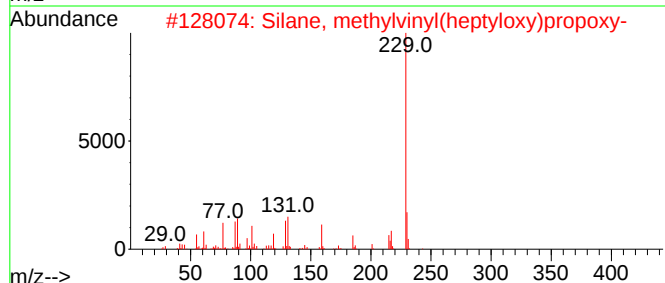
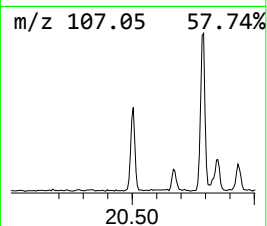
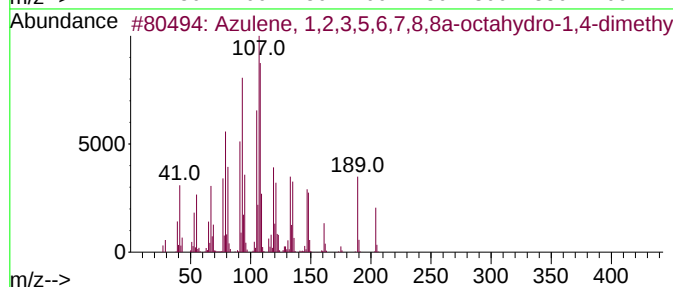
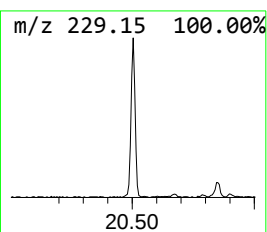
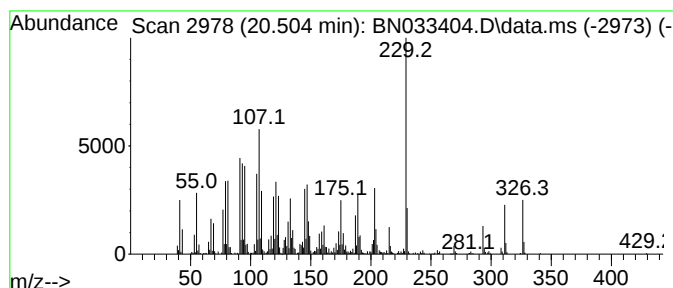
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 unknown-02 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.504	2.52 ng/ul	727708	Chrysene-d12	21.192

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Azulene, 1,2,3,5,6,7,8,8a-octahy...	204	C15H24	003691-11-0	25
2			Silane, methylvinyl(heptyloxy)pr...	244	C13H28O2Si	1000416-86-2	22
3			Celestolide	244	C17H24O	013171-00-1	22
4			Silane, diethyl(cis-4-methylcycl...	258	C14H30O2Si	1000363-54-8	22
5			1-Piperazinepropanenitrile, .bet...	229	C13H15N3O	014761-40-1	22



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081524\
Data File : BN033404.D
Acq On : 15 Aug 2024 17:35
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Sample : P3481-19
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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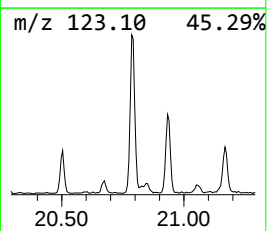
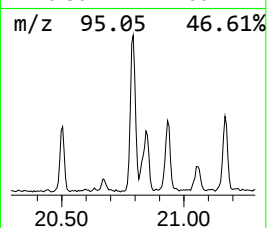
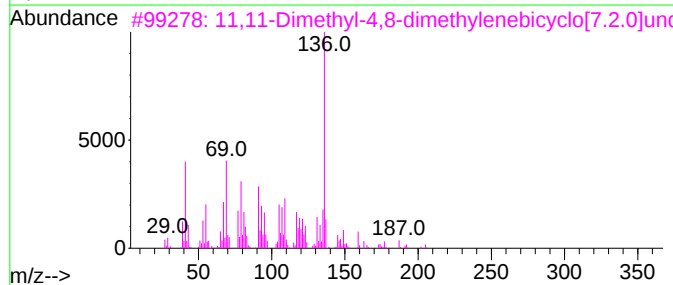
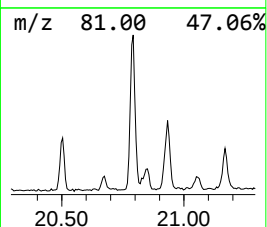
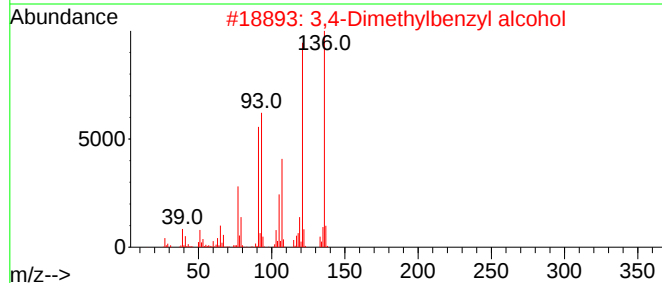
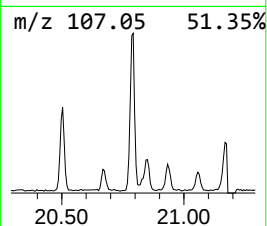
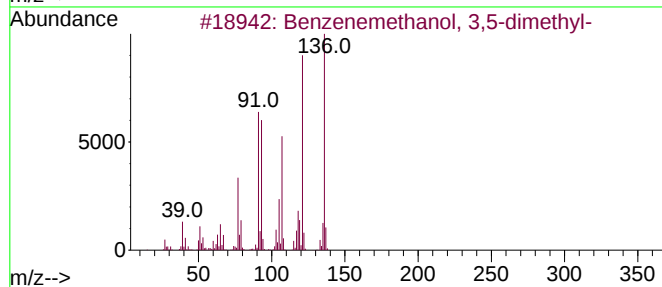
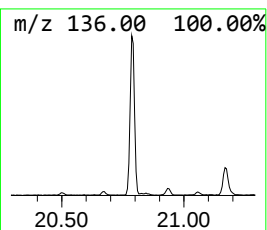
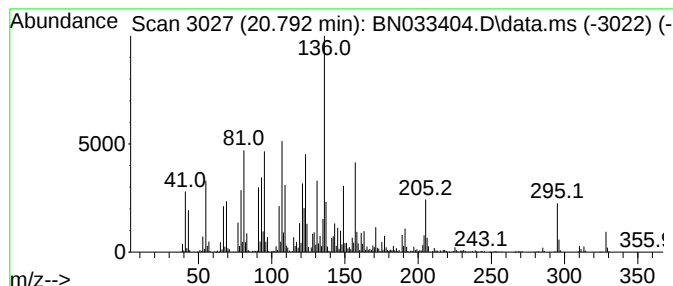
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 unknown-03 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.792	6.06 ng/ul	1749540	Chrysene-d12	21.192

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzenemethanol, 3,5-dimethyl-	136	C9H12O	027129-87-9	30
2			3,4-Dimethylbenzyl alcohol	136	C9H12O	006966-10-5	22
3			11,11-Dimethyl-4,8-dimethylenebi...	220	C15H24O	079580-01-1	22
4			o-Toluic acid, 2-tridecyl ester	318	C21H34O2	1000299-79-0	14
5			Phenol, 3-(ethylamino)-4-methyl-	151	C9H13NO	000120-37-6	14



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081524\
Data File : BN033404.D
Acq On : 15 Aug 2024 17:35
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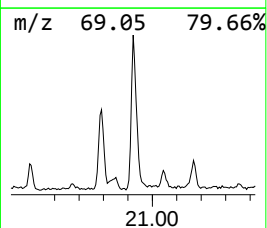
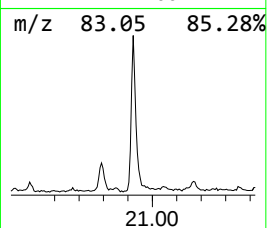
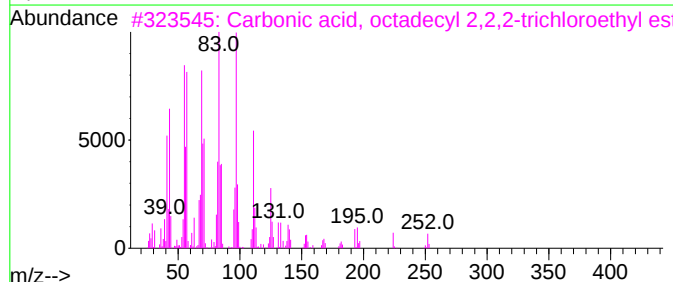
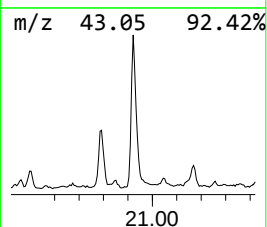
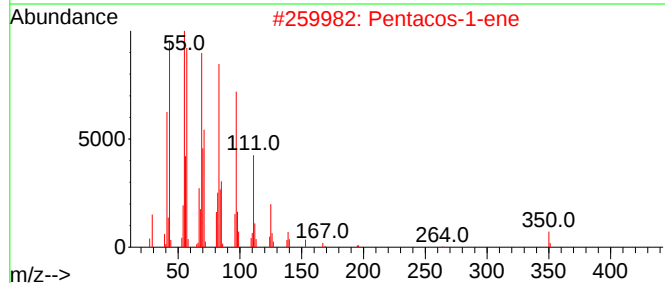
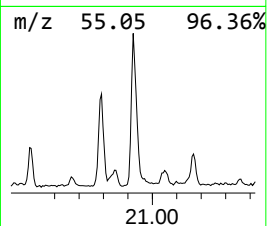
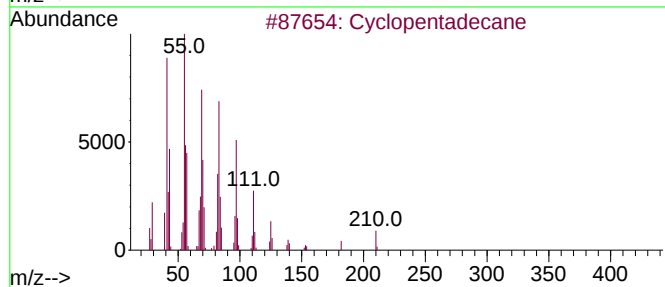
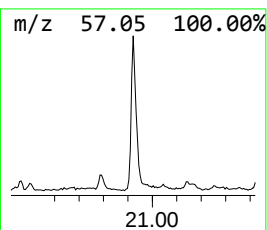
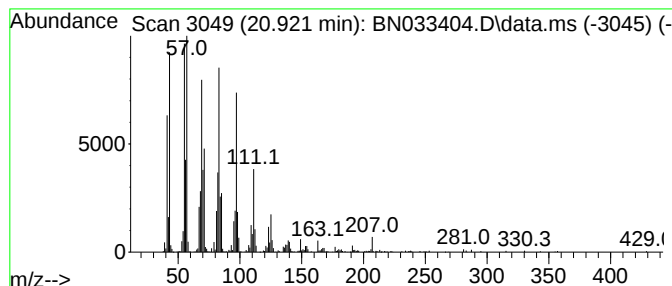
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 (DEL) Alkane: Cyclic20.921 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.921	4.12 ng/ul	1189600	Chrysene-d12	21.192

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopentadecane	210	C15H30	000295-48-7	97
2			Pentacos-1-ene	350	C25H50	016980-85-1	94
3			Carbonic acid, octadecyl 2,2,2-t...	444	C21H39Cl3O3	1000314-56-3	94
4			n-Tetracosanol-1	354	C24H50O	000506-51-4	94
5			Trifluoroacetoxy hexadecane	338	C18H33F3O2	006222-03-3	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081524\
Data File : BN033404.D
Acq On : 15 Aug 2024 17:35
Operator : MA/JU
Sample : P3481-19
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST0.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
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1-Phenoxypropan...	11.081	3.0	ng/ul	617085	2	10.334	4163030	20.0
unknown-01	14.463	2.9	ng/ul	682427	3	14.210	4676770	20.0
Phenol, 4-(1-me...	17.345	16.5	ng/ul	3970610	4	16.957	4822080	20.0
n-Hexadecanoic ...	17.892	2.8	ng/ul	661944	4	16.957	4822080	20.0
Phenol, 4,4'-(1...	19.421	9.6	ng/ul	2780160	5	21.192	5775000	20.0
unknown-02	20.504	2.5	ng/ul	727708	5	21.192	5775000	20.0
unknown-03	20.792	6.1	ng/ul	1749540	5	21.192	5775000	20.0
(DEL) Alkane: C...	20.921	4.1	ng/ul	1189600	5	21.192	5775000	20.0