

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081524\
 Data File : BN033403.D
 Acq On : 15 Aug 2024 16:56
 Operator : MA/JU
 Sample : P3481-18
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 E0AF3

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: BN033403.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.158	25	29	39	rVB	146684	203890	1.70%	0.127%
2	3.411	68	72	88	rBV	577982	783755	6.53%	0.488%
3	3.546	91	95	111	rVV	1446627	2085219	17.37%	1.297%
4	3.864	145	149	156	rVB	30962	45594	0.38%	0.028%
5	4.499	254	257	267	rVB	21309	39365	0.33%	0.024%
6	4.746	293	299	305	rVB	10115	18233	0.15%	0.011%
7	4.975	330	338	345	rBV	38344	65648	0.55%	0.041%
8	5.322	392	397	406	rVB2	12358	19201	0.16%	0.012%
9	5.693	455	460	467	rVB	84953	119879	1.00%	0.075%
10	6.516	595	600	606	rVB2	38201	53902	0.45%	0.034%
11	6.587	606	612	618	rBV2	34728	57147	0.48%	0.036%
12	6.752	629	640	656	rVB	3010417	4621796	38.49%	2.876%
13	6.875	656	661	662	rVV	18121	24550	0.20%	0.015%
14	6.916	662	668	680	rVB	2509888	3691161	30.74%	2.297%
15	7.105	691	700	708	rBV	2947343	4635669	38.61%	2.884%
16	7.181	708	713	726	rVB	348564	517701	4.31%	0.322%
17	7.569	771	779	786	rBV	2279130	3527608	29.38%	2.195%
18	7.646	786	792	798	rVB	68525	99532	0.83%	0.062%
19	7.875	826	831	836	rBV3	14376	25213	0.21%	0.016%
20	8.275	891	899	911	rBV	3659442	5748384	47.88%	3.576%
21	8.710	965	973	983	rBV	2838763	4477233	37.29%	2.786%
22	8.846	991	996	1000	rVV3	15559	23158	0.19%	0.014%
23	8.893	1000	1004	1010	rVB6	13577	21287	0.18%	0.013%
24	9.193	1045	1055	1061	rVB	24805	49540	0.41%	0.031%
25	9.287	1064	1071	1079	rVV	284409	423914	3.53%	0.264%
26	9.428	1085	1095	1103	rBV	2170397	3541825	29.50%	2.204%
27	9.663	1130	1135	1142	rVB4	13740	24761	0.21%	0.015%
28	9.957	1178	1185	1196	rVV	3566071	5750090	47.89%	3.578%
29	10.134	1208	1215	1223	rVB4	19459	39544	0.33%	0.025%
30	10.334	1240	1249	1258	rBV	3120359	5192042	43.24%	3.230%
31	10.469	1264	1272	1282	rBV	1504160	2454257	20.44%	1.527%
32	10.893	1339	1344	1350	rBV	126761	184721	1.54%	0.115%
33	11.081	1369	1376	1383	rBV	704746	1128061	9.40%	0.702%
34	11.140	1383	1386	1394	rVB5	19747	34792	0.29%	0.022%
35	11.799	1491	1498	1502	rBV8	8760	17575	0.15%	0.011%
36	11.846	1502	1506	1513	rVV2	10139	18139	0.15%	0.011%

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Integrator: RTE

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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

37	11.928	1514	1520	1527	rVB	79462	132164	1.10%	0.082%
38	12.493	1607	1616	1619	rBV10	8132	18236	0.15%	0.011%
39	12.569	1625	1629	1633	rBV2	13241	20895	0.17%	0.013%
40	12.822	1667	1672	1676	rBV2	18596	28322	0.24%	0.018%

41	12.987	1692	1700	1707	rBV5	15321	30149	0.25%	0.019%
42	13.063	1707	1713	1718	rBV2	26791	35523	0.30%	0.022%
43	13.540	1791	1794	1799	rVB5	36739	51496	0.43%	0.032%
44	13.634	1803	1810	1816	rVV	5589532	7867711	65.53%	4.895%
45	13.898	1847	1855	1866	rBV	6811329	10101086	84.13%	6.285%

46	14.069	1876	1884	1891	rBV8	34667	64840	0.54%	0.040%
47	14.210	1901	1908	1916	rBV	4544859	6631683	55.23%	4.126%
48	14.287	1916	1921	1927	rVB8	42643	65021	0.54%	0.040%
49	14.345	1927	1931	1935	rBV3	43285	60971	0.51%	0.038%
50	14.410	1935	1942	1946	rBV	2443744	3697046	30.79%	2.300%

51	14.445	1946	1948	1950	rBV	332311	397841	3.31%	0.248%
52	14.657	1977	1984	1987	rBV7	23422	45076	0.38%	0.028%
53	14.740	1994	1998	2010	rVV2	42852	93153	0.78%	0.058%
54	14.881	2020	2022	2031	rVB10	9315	18486	0.15%	0.012%
55	15.045	2046	2050	2054	rBV4	33064	51537	0.43%	0.032%

56	15.110	2058	2061	2064	rVB	17969	19540	0.16%	0.012%
57	15.210	2071	2078	2091	rVB	8003542	11663759	97.15%	7.257%
58	15.322	2091	2097	2104	rBV	2059594	2793520	23.27%	1.738%
59	15.445	2114	2118	2128	rVB3	39530	64254	0.54%	0.040%
60	15.669	2152	2156	2160	rVB6	15304	19882	0.17%	0.012%

61	15.787	2170	2176	2184	rBV2	83167	131120	1.09%	0.082%
62	15.969	2204	2207	2210	rBV	32275	44619	0.37%	0.028%
63	16.128	2231	2234	2242	rVB7	21277	36262	0.30%	0.023%
64	16.392	2274	2279	2284	rBV2	54840	73443	0.61%	0.046%
65	16.551	2303	2306	2310	rBV4	21067	28923	0.24%	0.018%

66	16.716	2328	2334	2336	rBV5	21954	36646	0.31%	0.023%
67	16.763	2340	2342	2346	rVB2	20733	26612	0.22%	0.017%
68	16.957	2366	2375	2381	rVV	5271207	7452711	62.07%	4.637%
69	16.998	2381	2382	2385	rVV	56765	48275	0.40%	0.030%
70	17.057	2385	2392	2403	rVB	8446564	11795765	98.25%	7.339%

71	17.163	2405	2410	2417	rBV5	69964	95668	0.80%	0.060%
72	17.345	2435	2441	2463	rBV	1096234	2312262	19.26%	1.439%
73	17.675	2493	2497	2502	rVB	167419	199409	1.66%	0.124%
74	17.763	2507	2512	2519	rBV3	101263	200307	1.67%	0.125%
75	17.898	2528	2535	2551	rBV	1555543	2466797	20.55%	1.535%

76	18.181	2579	2583	2590	rVB3	57300	109069	0.91%	0.068%
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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	18.333	2605	2609	2616	rBV8	35948	67398	0.56%	0.042%
78	18.563	2645	2648	2656	rBV6	23581	41738	0.35%	0.026%
79	18.763	2678	2682	2688	rBV5	55608	89143	0.74%	0.055%
80	18.857	2694	2698	2702	rVB2	162586	199043	1.66%	0.124%
81	18.922	2705	2709	2715	rVB	123292	168091	1.40%	0.105%
82	18.992	2716	2721	2723	rBV	153339	205177	1.71%	0.128%
83	19.028	2723	2727	2730	rVV	136536	214525	1.79%	0.133%
84	19.057	2730	2732	2744	rVB5	86981	174626	1.45%	0.109%
85	19.186	2749	2754	2774	rBV	705206	1098772	9.15%	0.684%
86	19.363	2777	2784	2790	rBV	8600421	12006317	100.00%	7.470%
87	19.422	2790	2794	2817	rVB	1267602	2129490	17.74%	1.325%
88	19.922	2876	2879	2883	rVB4	42314	43174	0.36%	0.027%
89	19.963	2884	2886	2889	rVB	34311	31460	0.26%	0.020%
90	20.922	3045	3049	3065	rBV	1217493	1913827	15.94%	1.191%
91	21.110	3078	3081	3084	rBV	81460	90936	0.76%	0.057%
92	21.192	3089	3095	3101	rVV	3930350	5388123	44.88%	3.352%
93	21.439	3133	3137	3141	rVB2	99484	133399	1.11%	0.083%
94	21.986	3225	3230	3242	rVB	566693	961054	8.00%	0.598%
95	22.604	3332	3335	3342	rVB3	84672	138835	1.16%	0.086%
96	22.763	3358	3362	3368	rVB	222090	351236	2.93%	0.219%
97	23.316	3450	3456	3477	rVB	584040	1312685	10.93%	0.817%
98	23.833	3534	3544	3566	rVB	3953114	9436103	78.59%	5.871%
99	24.021	3567	3576	3586	rBV	2271966	5254884	43.77%	3.269%
100	25.221	3775	3780	3796	rVB5	166268	482234	4.02%	0.300%

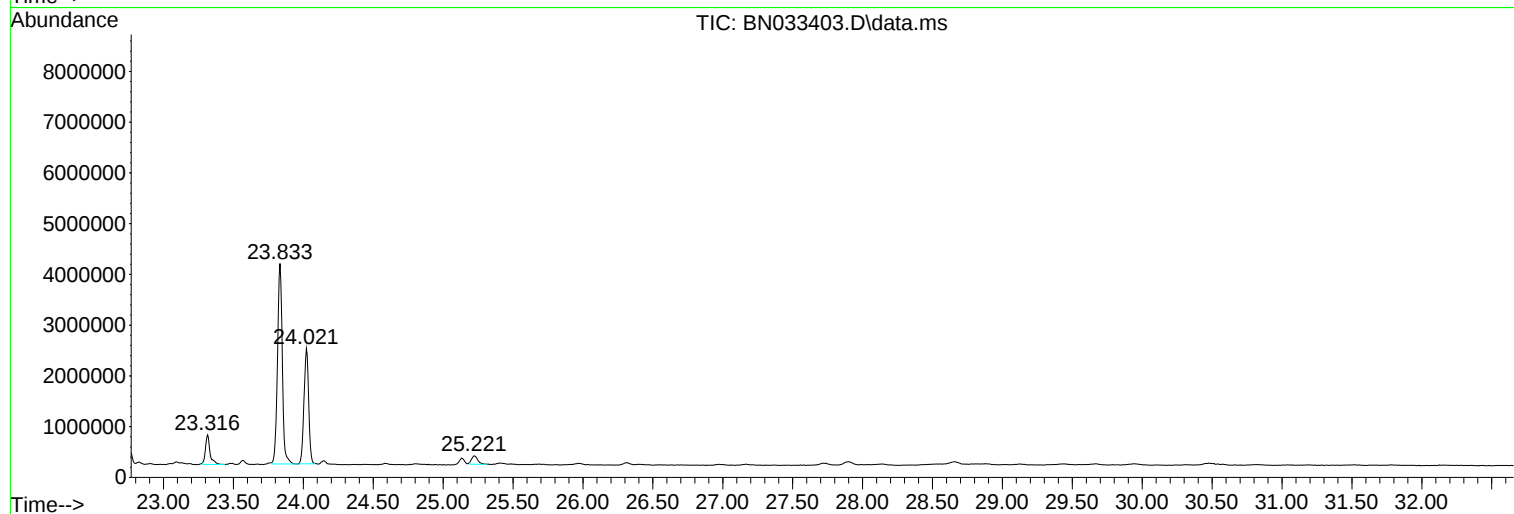
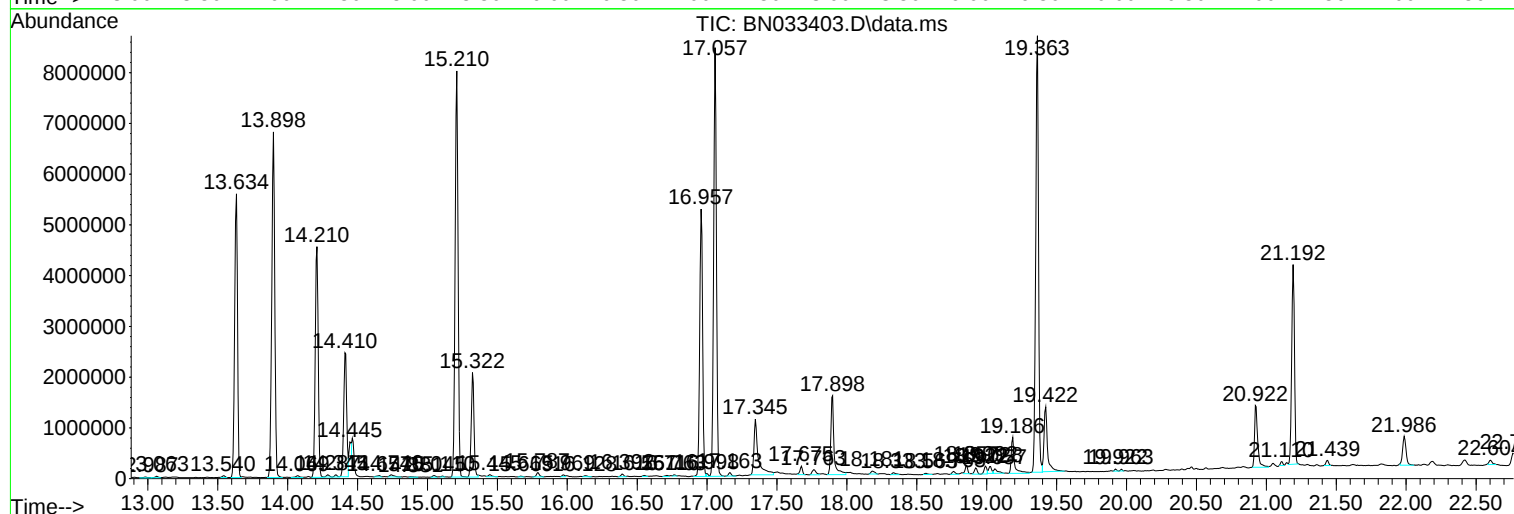
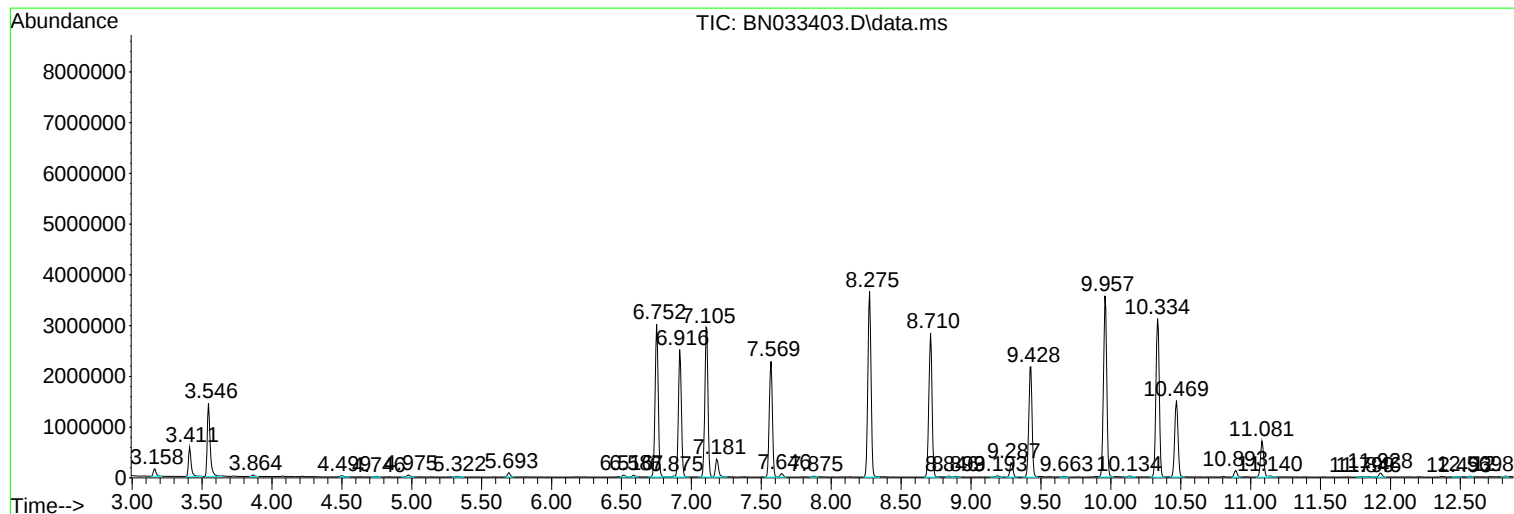
Sum of corrected areas: 160726735

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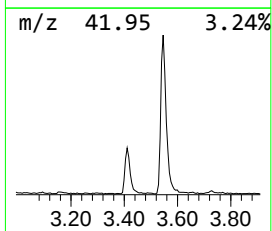
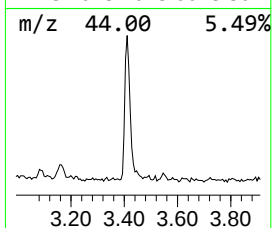
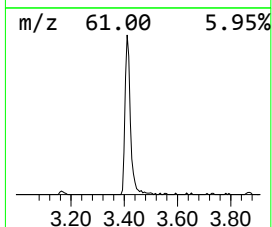
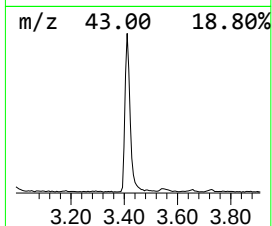
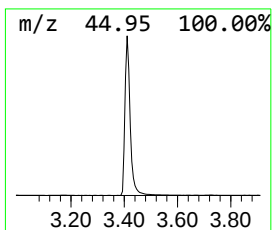
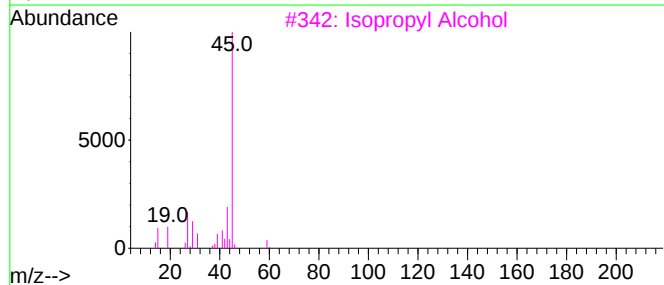
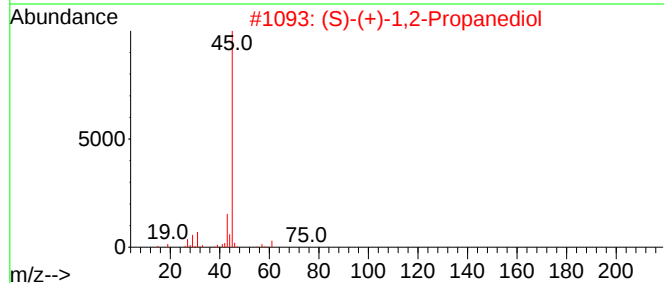
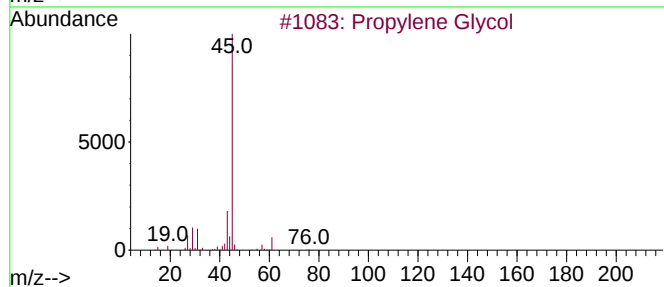
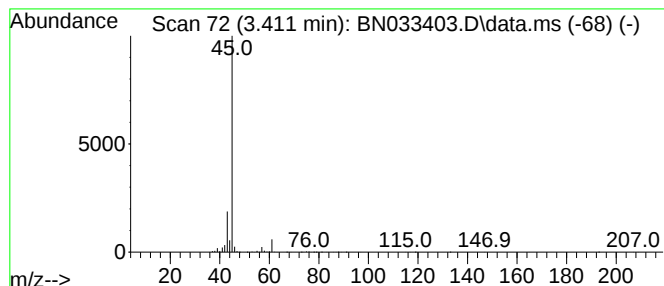
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN081424.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Propylene Glycol Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.411	4.44 ng/ul	783755	1,4-Dichlorobenzene-d4	7.569

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Propylene Glycol	76	C3H8O2	000057-55-6	86
2			(S)-(+)-1,2-Propanediol	76	C3H8O2	004254-15-3	64
3			Isopropyl Alcohol	60	C3H8O	000067-63-0	56
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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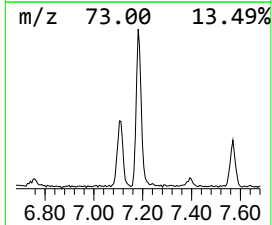
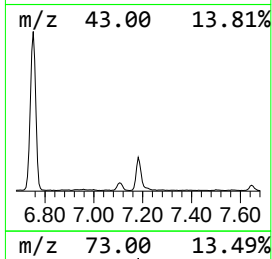
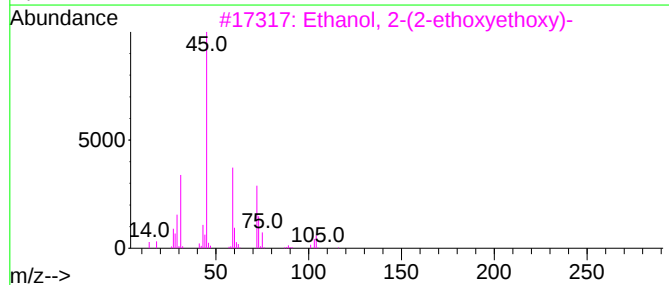
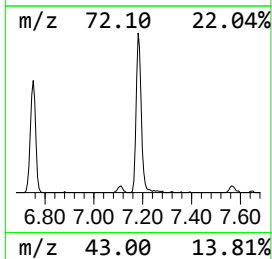
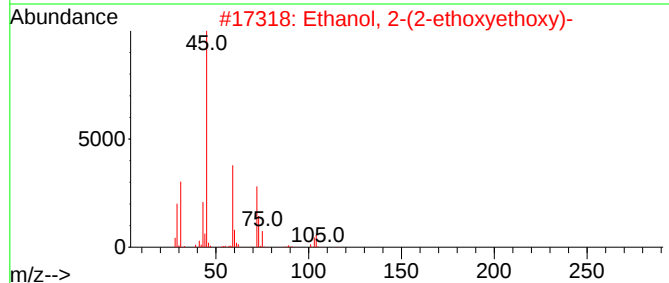
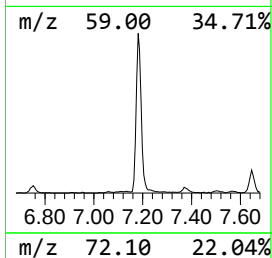
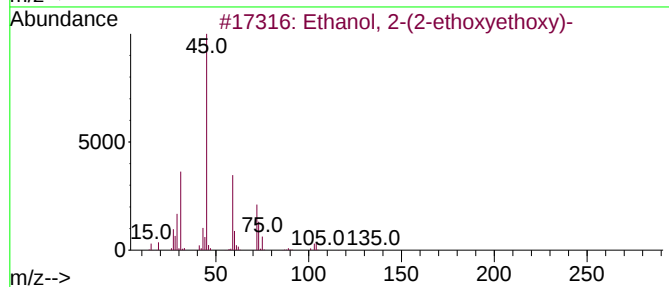
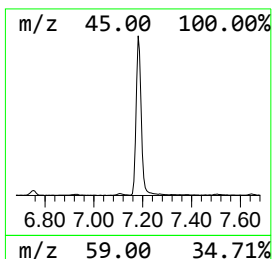
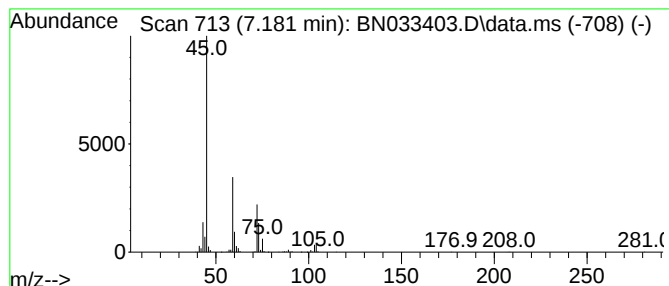
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN081424.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Ethanol, 2-(2-ethoxyethoxy)- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.181	2.94 ng/ul	517701	1,4-Dichlorobenzene-d4	7.569

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol, 2-(2-ethoxyethoxy)-	134	C6H14O3	000111-90-0	90
2			Ethanol, 2-(2-ethoxyethoxy)-	134	C6H14O3	000111-90-0	90
3			Ethanol, 2-(2-ethoxyethoxy)-	134	C6H14O3	000111-90-0	78
4			Diethyl carbitol	162	C8H18O3	000112-36-7	72
5			Ethanol, 2-[2-(2-ethoxyethoxy)et...	178	C8H18O4	000112-50-5	64



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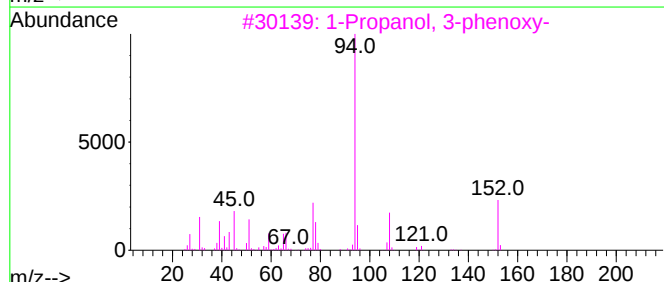
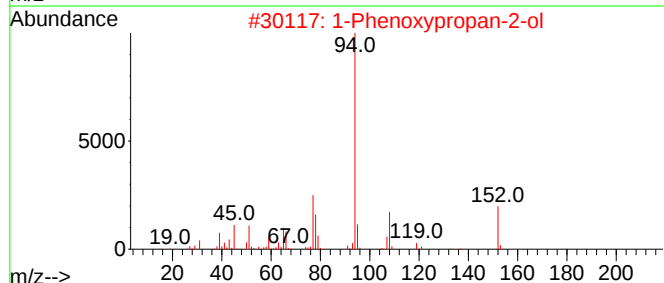
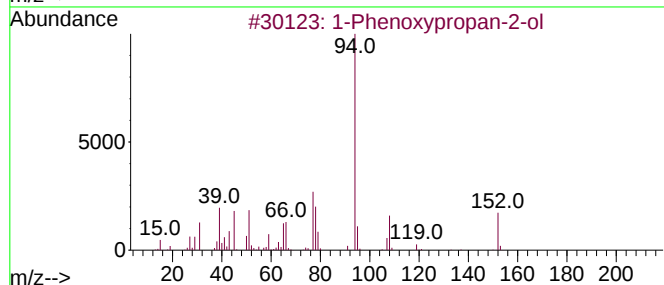
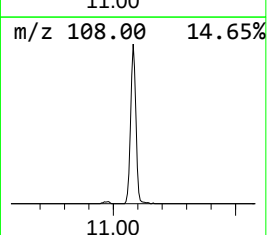
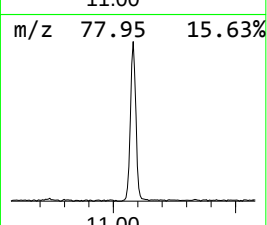
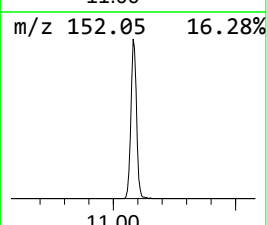
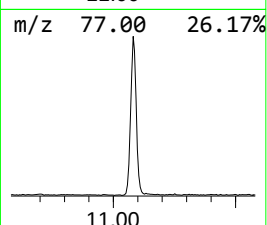
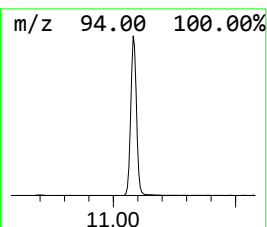
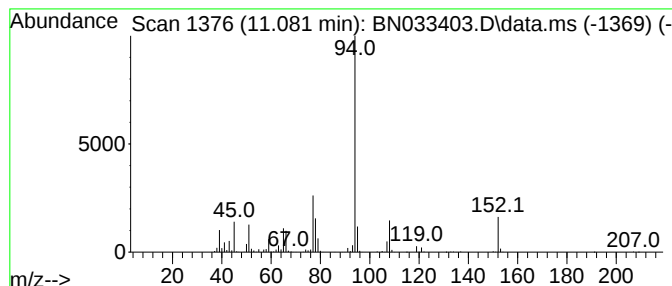
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.081	4.35 ng/ul	1128060	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	96
3			1-Propanol, 3-phenoxy-	152	C9H12O2	006180-61-6	93
4			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	91
5			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081524\
Data File : BN033403.D
Acq On : 15 Aug 2024 16:56
Operator : MA/JU
Sample : P3481-18
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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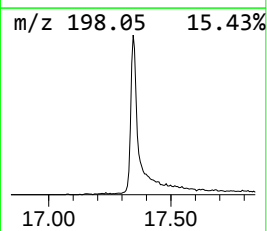
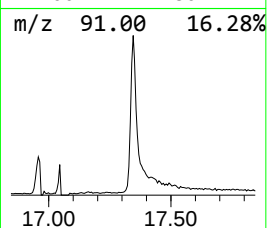
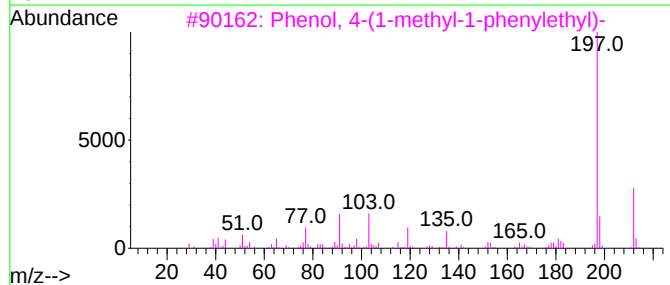
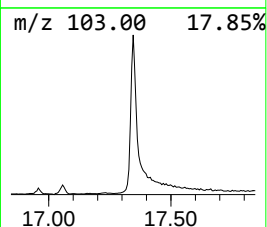
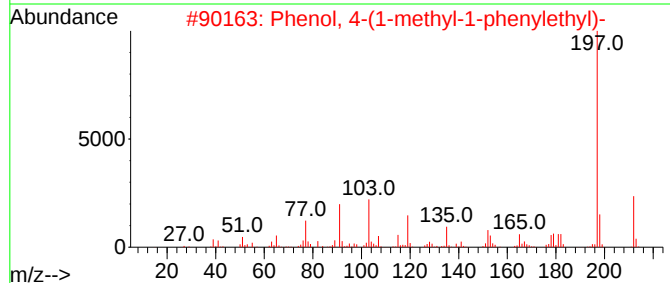
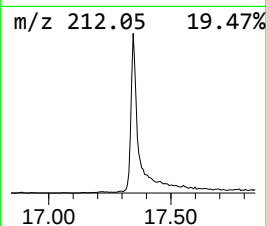
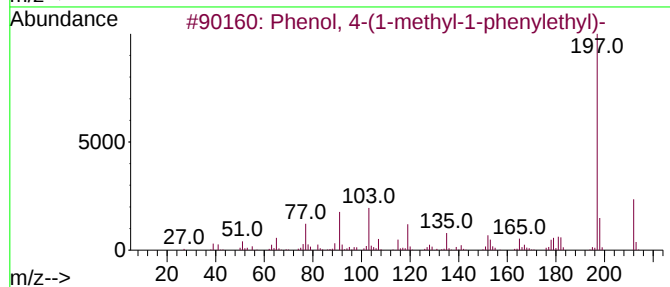
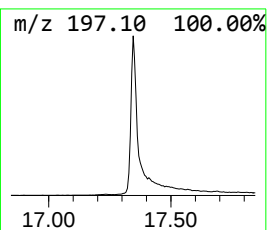
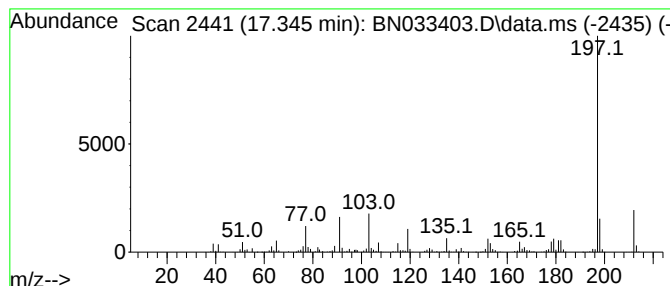
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.345	6.21 ng/ul	2312260	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	99
2			Phenol, 4-(1-methyl-1-phenylethyl)-	212	C15H16O	000599-64-4	98
3			Phenol, 4-(1-methyl-1-phenylethyl)-	212	C15H16O	000599-64-4	87
4			Dimethyl diphenylsilane	212	C14H16Si	000778-24-5	68
5			2,3-Dimethyl-4-biphenylamine	197	C14H15N	022750-87-4	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081524\
Data File : BN033403.D
Acq On : 15 Aug 2024 16:56
Operator : MA/JU
Sample : P3481-18
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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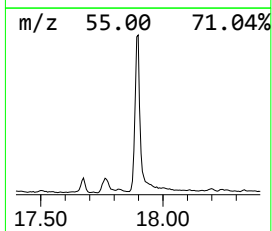
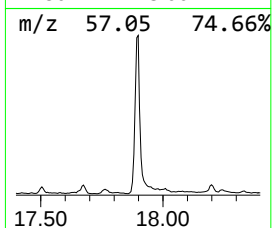
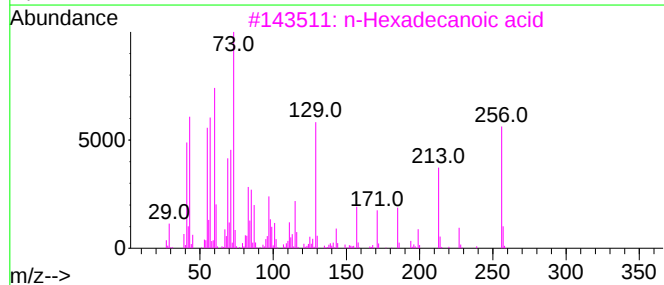
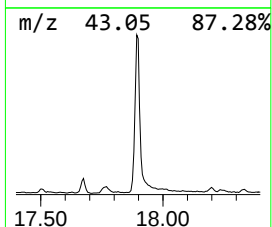
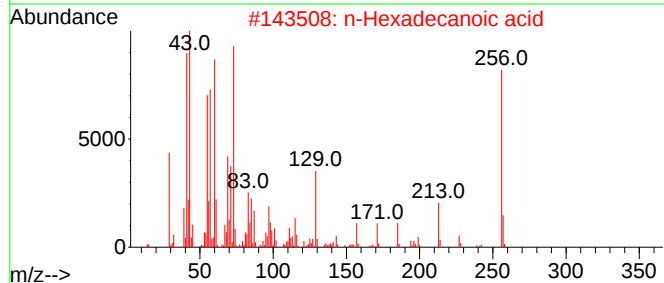
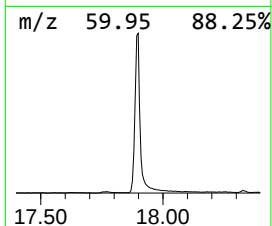
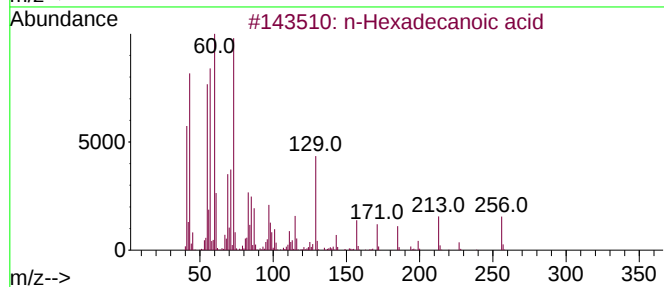
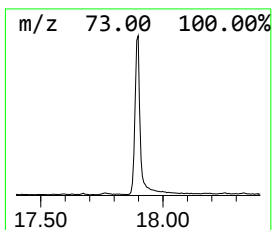
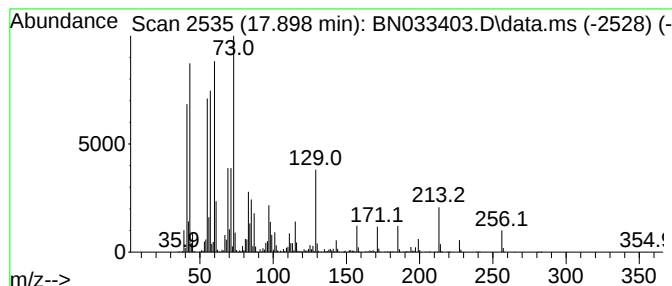
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.898	6.62 ng/ul	2466800	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	99
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
5			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081524\
Data File : BN033403.D
Acq On : 15 Aug 2024 16:56
Operator : MA/JU
Sample : P3481-18
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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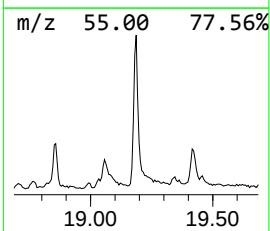
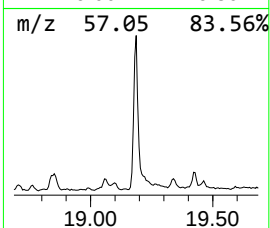
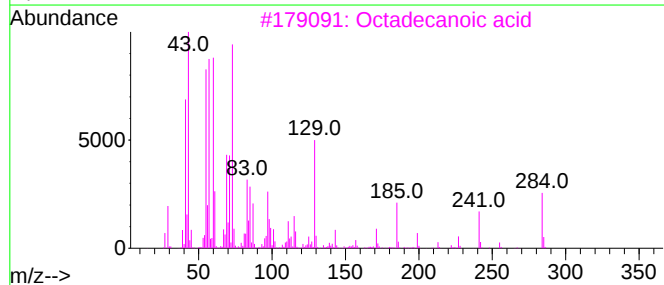
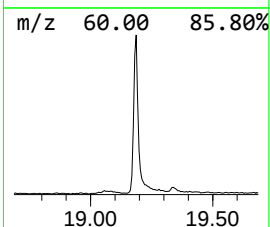
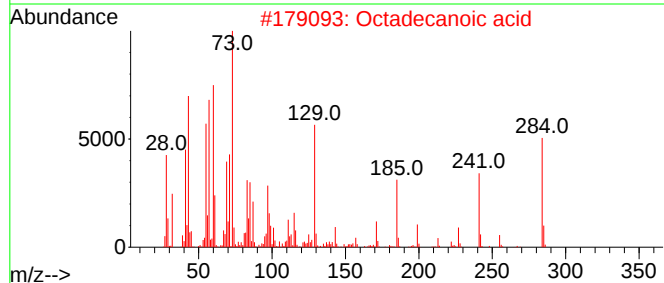
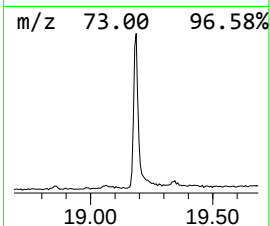
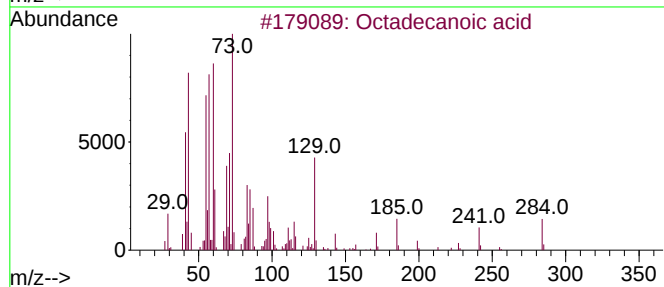
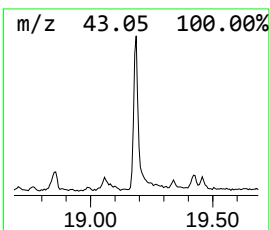
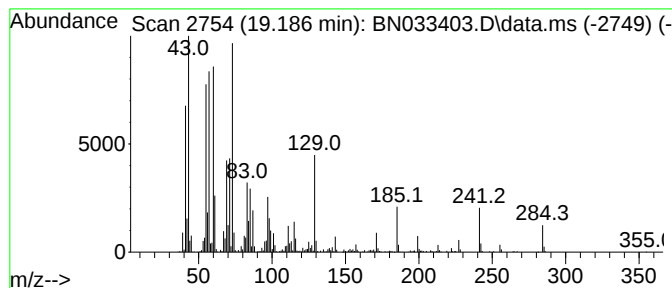
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Octadecanoic acid Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.186	4.08 ng/ul	1098770	Chrysene-d12	21.192

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid	284	C18H36O2	000057-11-4	99
2			Octadecanoic acid	284	C18H36O2	000057-11-4	99
3			Octadecanoic acid	284	C18H36O2	000057-11-4	99
4			Octadecanoic acid	284	C18H36O2	000057-11-4	95
5			Octadecanoic acid	284	C18H36O2	000057-11-4	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081524\
Data File : BN033403.D
Acq On : 15 Aug 2024 16:56
Operator : MA/JU
Sample : P3481-18
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
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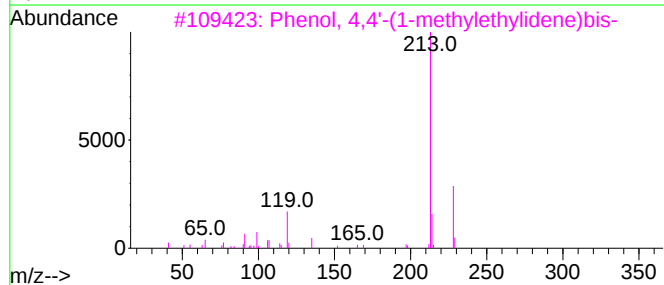
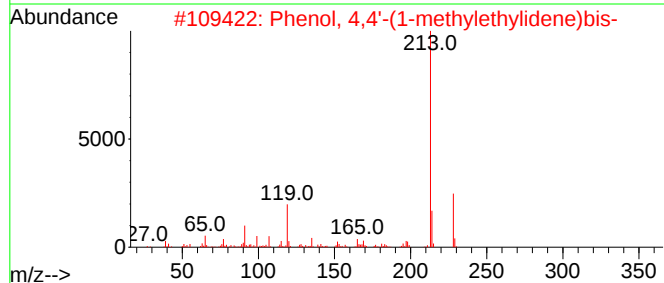
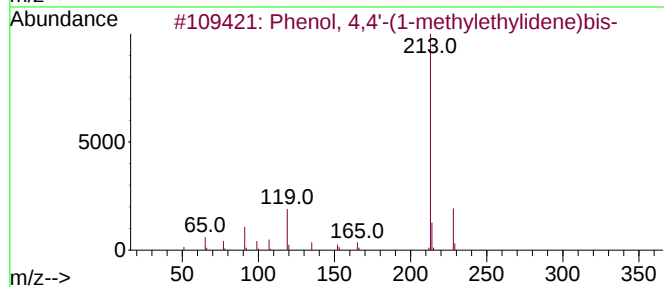
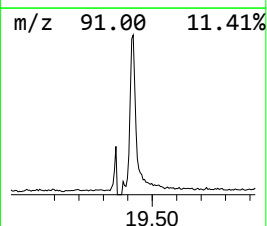
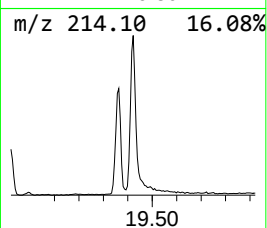
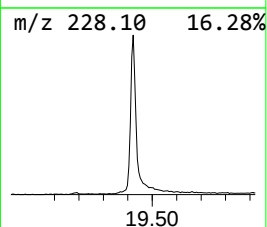
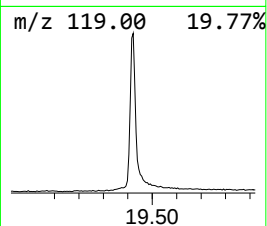
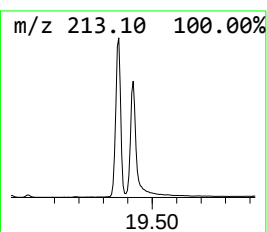
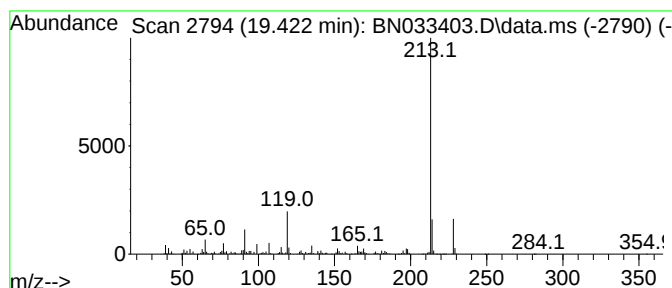
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Phenol, 4,4'-(1-methylethyl... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.422	7.90 ng/ul	2129490	Chrysene-d12	21.192

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenol, 4,4'-(1-methylethylidene...	228	C15H16O2	000080-05-7	97
2			Phenol, 4,4'-(1-methylethylidene...	228	C15H16O2	000080-05-7	95
3			Phenol, 4,4'-(1-methylethylidene...	228	C15H16O2	000080-05-7	91
4			3,4'-Isopropylidenediphenol	228	C15H16O2	046765-25-7	78
5			4,4'-(1,3-Dimethylbutylidene)bis...	270	C18H22O2	006807-17-6	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081524\
Data File : BN033403.D
Acq On : 15 Aug 2024 16:56
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Misc :
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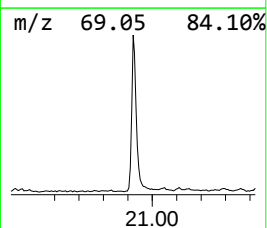
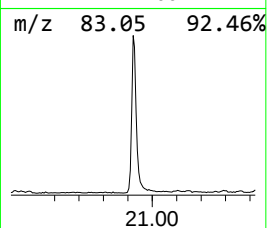
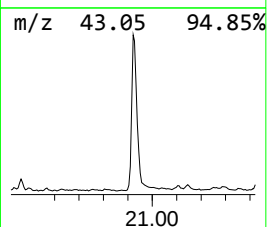
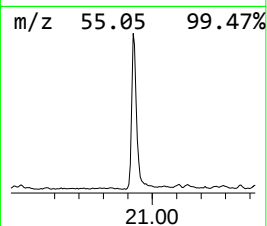
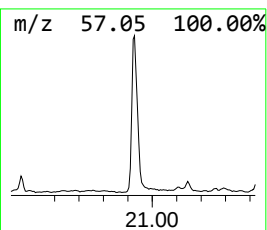
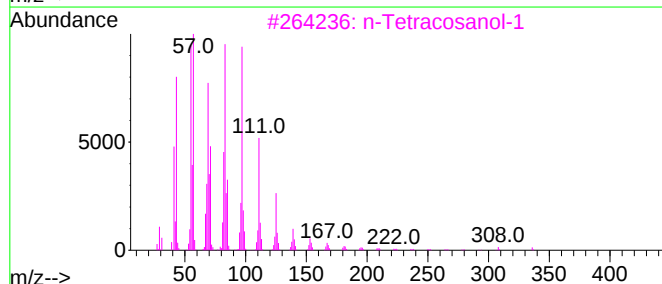
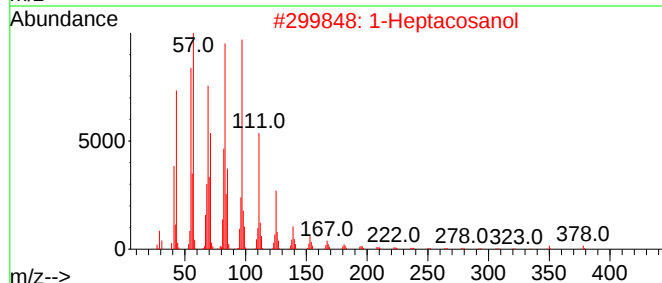
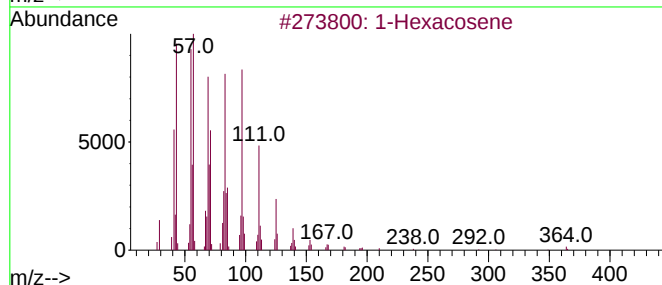
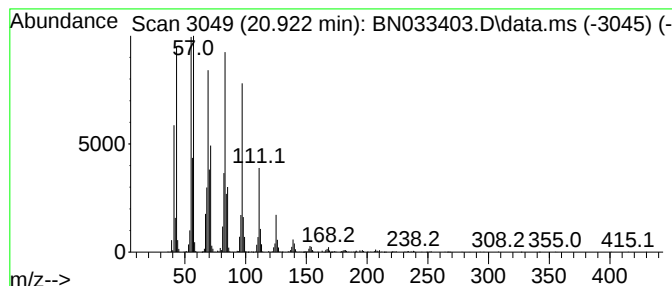
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.922	7.10 ng/ul	1913830	Chrysene-d12	21.192

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Hexacosene			364	C26H52	018835-33-1	94
2	1-Heptacosanol			396	C27H56O	002004-39-9	94
3	n-Tetracosanol-1			354	C24H50O	000506-51-4	94
4	Behenic alcohol			326	C22H46O	000661-19-8	94
5	Trichloroacetic acid, hexadecyl ...			386	C18H33Cl3O2	074339-54-1	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081524\
Data File : BN033403.D
Acq On : 15 Aug 2024 16:56
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ALS Vial : 12 Sample Multiplier: 1

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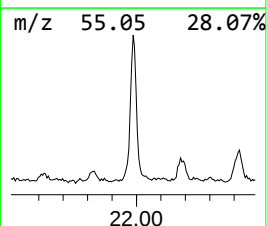
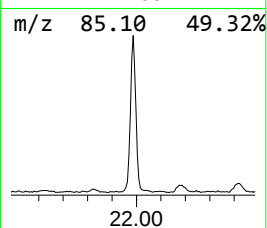
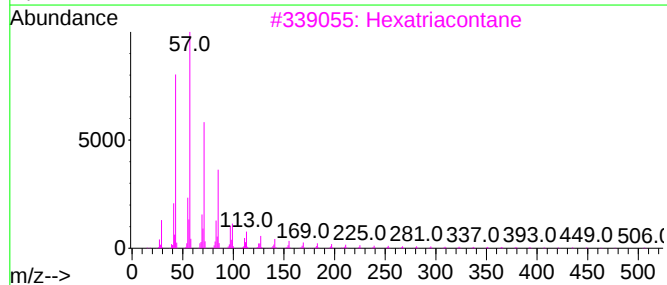
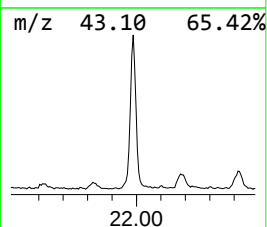
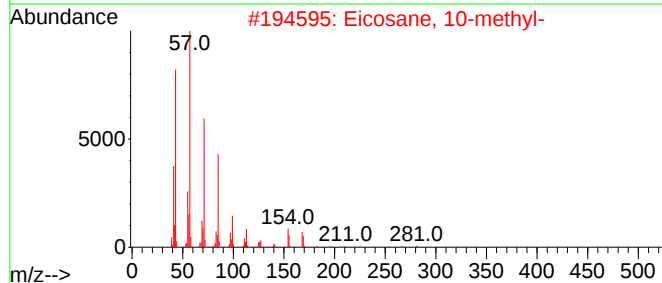
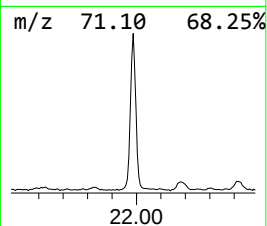
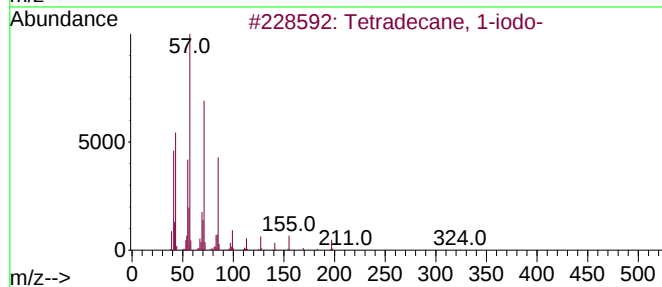
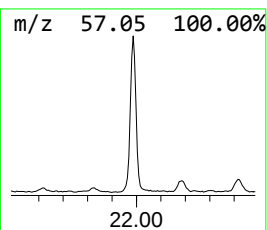
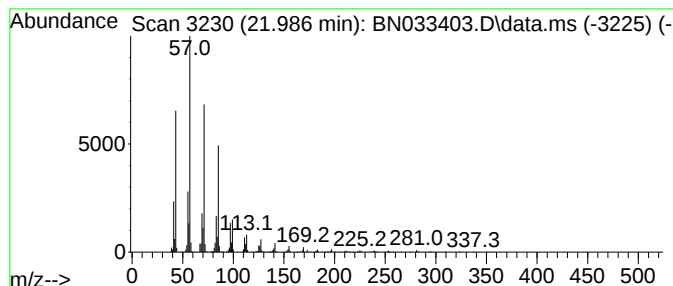
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 Tetradecane, 1-iodo- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.986	3.57 ng/ul	961054	Chrysene-d12	21.192

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Tetradecane, 1-iodo-	324	C14H29I	019218-94-1	91
2		Eicosane, 10-methyl-	296	C21H44	054833-23-7	91
3		Hexatriacontane	507	C36H74	000630-06-8	91
4		Hexacosane	366	C26H54	000630-01-3	91
5		Tetratetracontane	619	C44H90	007098-22-8	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081524\
Data File : BN033403.D
Acq On : 15 Aug 2024 16:56
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Sample : P3481-18
Misc :
ALS Vial : 12 Sample Multiplier: 1

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ClientSampleId :
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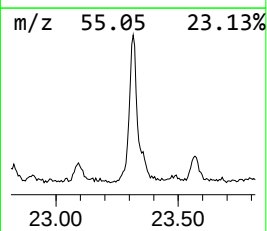
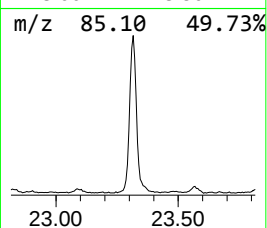
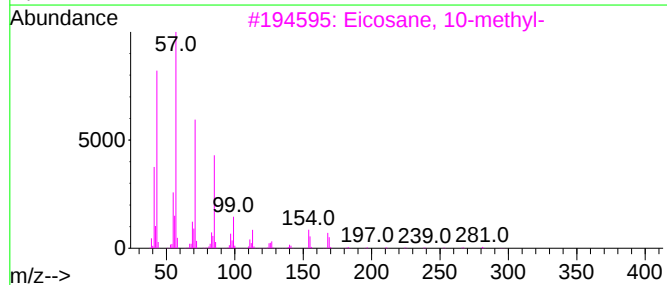
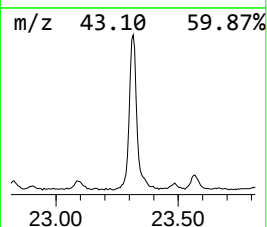
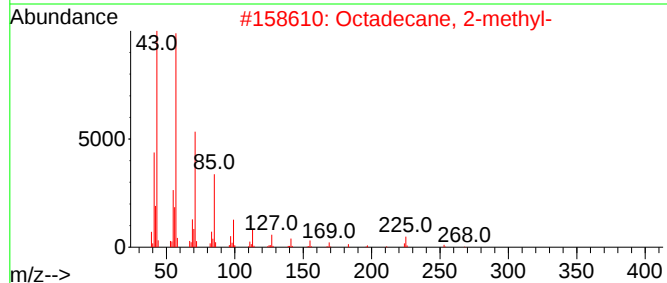
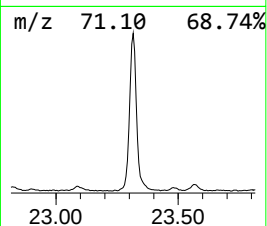
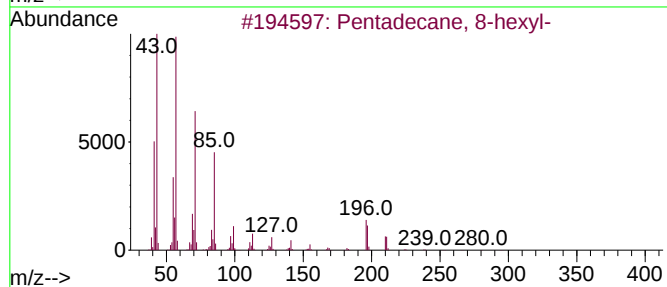
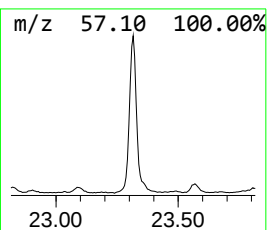
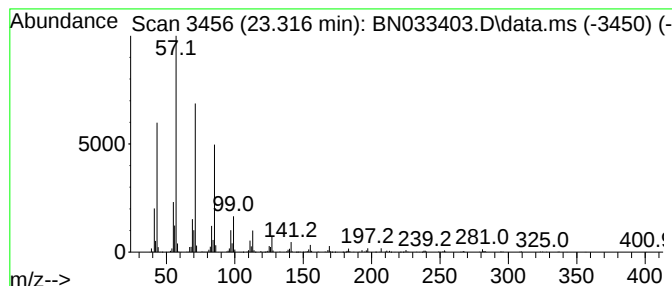
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.316	5.00 ng/ul	1312690	Perylene-d12	24.021

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	94
2		Octadecane, 2-methyl-	268	C19H40	001560-88-9	94
3		Eicosane, 10-methyl-	296	C21H44	054833-23-7	93
4		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
5		Hexacosane	366	C26H54	000630-01-3	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081524\
Data File : BN033403.D
Acq On : 15 Aug 2024 16:56
Operator : MA/JU
Sample : P3481-18
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
E0AF3

Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN081424.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
Propylene Glycol	3.411	4.4	ng/ul	783755	1	7.569	3527610	20.0
Ethanol, 2-(2-e...	7.181	2.9	ng/ul	517701	1	7.569	3527610	20.0
1-Phenoxypropan...	11.081	4.3	ng/ul	1128060	2	10.334	5192040	20.0
Phenol, 4-(1-me...	17.345	6.2	ng/ul	2312260	4	16.957	7452710	20.0
n-Hexadecanoic ...	17.898	6.6	ng/ul	2466800	4	16.957	7452710	20.0
Octadecanoic acid	19.186	4.1	ng/ul	1098770	5	21.192	5388120	20.0
Phenol, 4,4'-(1...	19.422	7.9	ng/ul	2129490	5	21.192	5388120	20.0
(DEL) Alkane: S...	20.922	7.1	ng/ul	1913830	5	21.192	5388120	20.0
Tetradecane, 1-...	21.986	3.6	ng/ul	961054	5	21.192	5388120	20.0
(DEL) Alkane: S...	23.316	5.0	ng/ul	1312690	6	24.021	5254880	20.0