

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN060524\
 Data File : BN031798.D
 Acq On : 05 Jun 2024 13:48
 Operator : MA/JU
 Sample : P2623-04
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BHBQ9

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: BN031798.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.187	29	34	48	rVB	51651	92049	1.18%	0.085%
2	3.446	74	78	90	rBV	307602	467515	5.99%	0.433%
3	3.593	100	103	118	rBV	619589	1077839	13.81%	0.998%
4	3.905	152	156	164	rBV	75375	117317	1.50%	0.109%
5	5.769	466	473	480	rVB	34932	58980	0.76%	0.055%
6	6.846	646	656	671	rVB	985751	1605220	20.57%	1.486%
7	7.016	678	685	695	rVB	798041	1282670	16.43%	1.188%
8	7.205	709	717	725	rBV	937458	1576073	20.19%	1.459%
9	7.275	725	729	743	rVB	147762	263524	3.38%	0.244%
10	7.675	789	797	804	rBV	931551	1514217	19.40%	1.402%
11	7.746	804	809	815	rVB2	28077	47898	0.61%	0.044%
12	8.381	908	917	926	rBV	1227535	2021158	25.90%	1.871%
13	8.828	985	993	1001	rBV	927521	1559572	19.98%	1.444%
14	9.393	1083	1089	1098	rVB	108089	171236	2.19%	0.159%
15	9.546	1105	1115	1127	rBV	761683	1289439	16.52%	1.194%
16	10.081	1198	1206	1223	rBV	1197651	2068086	26.50%	1.915%
17	10.457	1261	1270	1276	rBV	1360554	2343615	30.03%	2.170%
18	10.505	1276	1278	1283	rVB	42920	58819	0.75%	0.054%
19	10.599	1283	1294	1304	rBV	1243381	2123760	27.21%	1.966%
20	11.004	1359	1363	1371	rVB2	38483	64466	0.83%	0.060%
21	11.199	1388	1396	1411	rBV	310641	623508	7.99%	0.577%
22	12.051	1534	1541	1548	rBV2	25765	44128	0.57%	0.041%
23	12.122	1548	1553	1561	rVB	29740	51204	0.66%	0.047%
24	12.346	1586	1591	1599	rVB2	26627	44291	0.57%	0.041%
25	13.498	1777	1787	1790	rBV8	16247	44783	0.57%	0.041%
26	13.634	1805	1810	1816	rBV2	26230	45776	0.59%	0.042%
27	13.734	1821	1827	1842	rVB	2264185	3270238	41.90%	3.028%
28	13.969	1861	1867	1868	rBV	53742	69299	0.89%	0.064%
29	14.010	1868	1874	1887	rVB	2667046	4075187	52.22%	3.773%
30	14.316	1918	1926	1932	rBV	2270738	3396119	43.51%	3.144%
31	14.381	1932	1937	1945	rVV	105982	171551	2.20%	0.159%
32	14.528	1955	1962	1972	rVV	1199643	1948991	24.97%	1.804%
33	14.716	1989	1994	2003	rVV	97407	186813	2.39%	0.173%
34	15.169	2069	2071	2081	rVB4	21751	37602	0.48%	0.035%
35	15.310	2089	2095	2101	rBV	3299904	4962105	63.58%	4.594%
36	15.369	2101	2105	2110	rVB	103100	146907	1.88%	0.136%

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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	15.440	2111	2117	2124	rBV	1036313	1437259	18.42%	1.331%
38	15.545	2129	2135	2144	rVV7	24523	86319	1.11%	0.080%
39	15.663	2150	2155	2163	rVB	53685	96129	1.23%	0.089%
40	15.810	2175	2180	2186	rVV	47317	94145	1.21%	0.087%
41	15.869	2186	2190	2199	rVB6	27439	58157	0.75%	0.054%
42	16.040	2212	2219	2225	rVB5	36563	82611	1.06%	0.076%
43	16.351	2265	2272	2273	rBV5	23719	42982	0.55%	0.040%
44	16.375	2273	2276	2280	rVV2	31317	48857	0.63%	0.045%
45	16.722	2330	2335	2342	rVB	165610	250243	3.21%	0.232%
46	16.792	2342	2347	2350	rBV3	28695	48993	0.63%	0.045%
47	16.887	2358	2363	2372	rVB	94742	158608	2.03%	0.147%
48	17.063	2387	2393	2397	rBV	2872113	4262681	54.62%	3.946%
49	17.104	2397	2400	2405	rVB	1830095	2496786	31.99%	2.312%
50	17.163	2405	2410	2415	rBV	3708448	5297979	67.88%	4.905%
51	17.257	2421	2426	2431	rVB2	44963	77035	0.99%	0.071%
52	17.469	2454	2462	2475	rVV2	233771	432070	5.54%	0.400%
53	17.586	2476	2482	2487	rVV3	31187	69000	0.88%	0.064%
54	17.645	2488	2492	2500	rVB5	48086	86922	1.11%	0.080%
55	17.745	2504	2509	2513	rBV3	84879	111095	1.42%	0.103%
56	17.939	2533	2542	2543	rBV2	214233	266854	3.42%	0.247%
57	17.969	2543	2547	2557	rVV2	518995	1268910	16.26%	1.175%
58	18.069	2560	2564	2569	rVV	98601	143405	1.84%	0.133%
59	18.134	2569	2575	2579	rVV2	259186	474867	6.08%	0.440%
60	18.169	2579	2581	2588	rVB	131269	176089	2.26%	0.163%
61	18.257	2591	2596	2599	rBV5	35993	66183	0.85%	0.061%
62	18.292	2599	2602	2606	rVB2	44514	60528	0.78%	0.056%
63	18.445	2623	2628	2631	rBV	179306	232484	2.98%	0.215%
64	18.486	2631	2635	2641	rVB	201191	275215	3.53%	0.255%
65	18.692	2667	2670	2674	rVB2	50523	60180	0.77%	0.056%
66	18.739	2675	2678	2681	rVV	47725	59197	0.76%	0.055%
67	18.781	2682	2685	2690	rVB2	54403	70509	0.90%	0.065%
68	18.869	2695	2700	2704	rBV2	115723	200802	2.57%	0.186%
69	18.922	2706	2709	2713	rVB2	56414	76207	0.98%	0.071%
70	19.004	2719	2723	2730	rBV2	176210	349733	4.48%	0.324%
71	19.128	2736	2744	2751	rVB	3070771	4220325	54.07%	3.907%
72	19.228	2758	2761	2762	rBV2	54867	54588	0.70%	0.051%
73	19.257	2762	2766	2773	rVB2	161045	263656	3.38%	0.244%
74	19.375	2783	2786	2789	rVB	62756	73951	0.95%	0.068%
75	19.463	2795	2801	2804	rBV	5335340	7646259	97.97%	7.079%
76	19.492	2804	2806	2814	rVV	2548197	3055777	39.15%	2.829%

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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

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 Peak separation: 5

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77	19.563	2814	2818	2825	rVB2	118013	186121	2.38%	0.172%
78	19.675	2833	2837	2842	rBV	142210	178588	2.29%	0.165%
79	19.733	2844	2847	2854	rVB	145439	215054	2.76%	0.199%
80	19.816	2856	2861	2868	rBV5	167134	418633	5.36%	0.388%
81	19.880	2868	2872	2875	rBV	50585	63788	0.82%	0.059%
82	19.951	2879	2884	2887	rBV4	129713	257063	3.29%	0.238%
83	19.986	2887	2890	2894	rVB2	200515	284231	3.64%	0.263%
84	20.092	2905	2908	2911	rVB	71141	77161	0.99%	0.071%
85	20.145	2912	2917	2921	rVV2	200151	349945	4.48%	0.324%
86	20.292	2938	2942	2945	rBV	98399	126317	1.62%	0.117%
87	20.333	2946	2949	2956	rVV2	105325	179132	2.30%	0.166%
88	20.739	3014	3018	3024	rBV	448559	608748	7.80%	0.564%
89	20.910	3042	3047	3051	rBV3	228617	421584	5.40%	0.390%
90	20.951	3051	3054	3057	rVV	201827	262535	3.36%	0.243%
91	20.986	3057	3060	3067	rVV3	932181	1409676	18.06%	1.305%
92	21.051	3067	3071	3078	rVB	379662	602495	7.72%	0.558%
93	21.304	3106	3114	3118	rBV3	3991186	7804580	100.00%	7.226%
94	21.345	3118	3121	3129	rVB	1305373	1828628	23.43%	1.693%
95	22.669	3342	3346	3355	rVB3	352856	670683	8.59%	0.621%
96	23.316	3449	3456	3463	rBV	954223	2910830	37.30%	2.695%
97	23.968	3563	3567	3572	rBV	327776	673904	8.63%	0.624%
98	24.039	3572	3579	3586	rVV2	3143591	7393848	94.74%	6.845%
99	24.104	3586	3590	3602	rVB	789055	1721210	22.05%	1.594%
100	24.245	3605	3614	3622	rBV	2364172	6112248	78.32%	5.659%

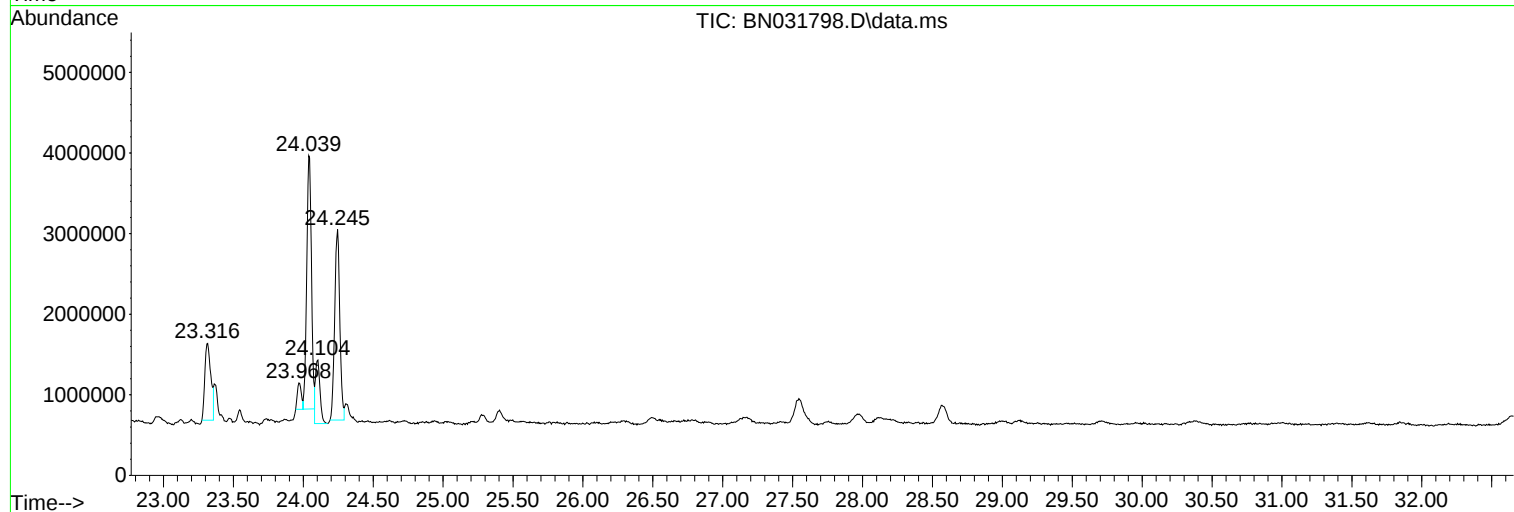
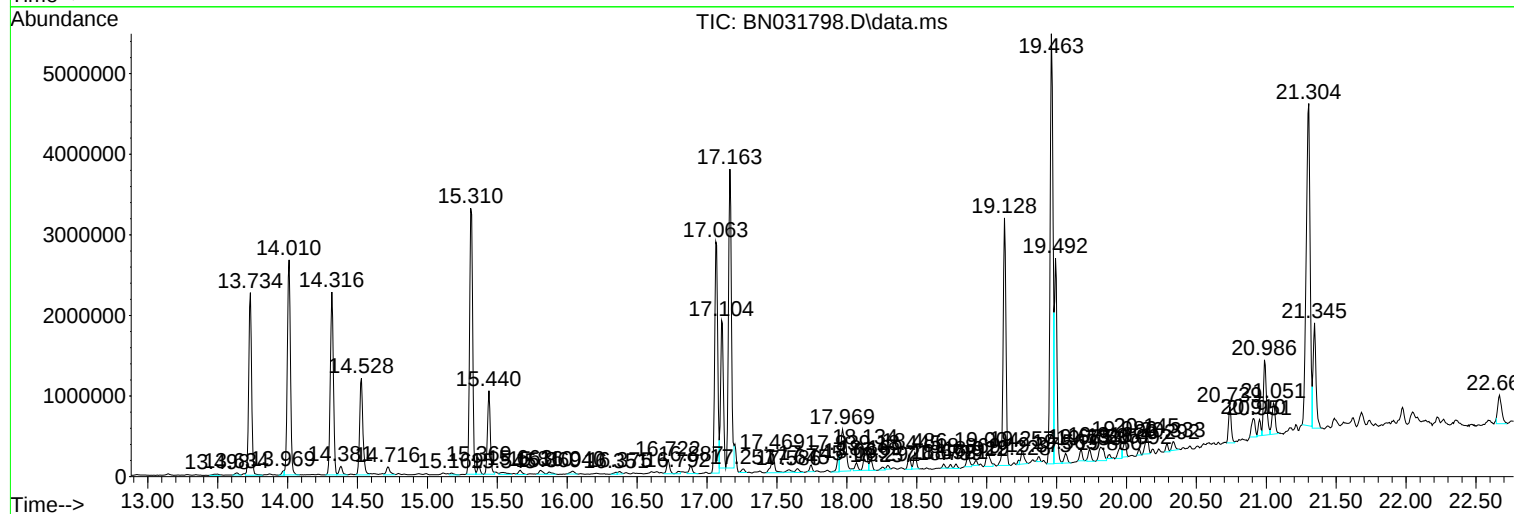
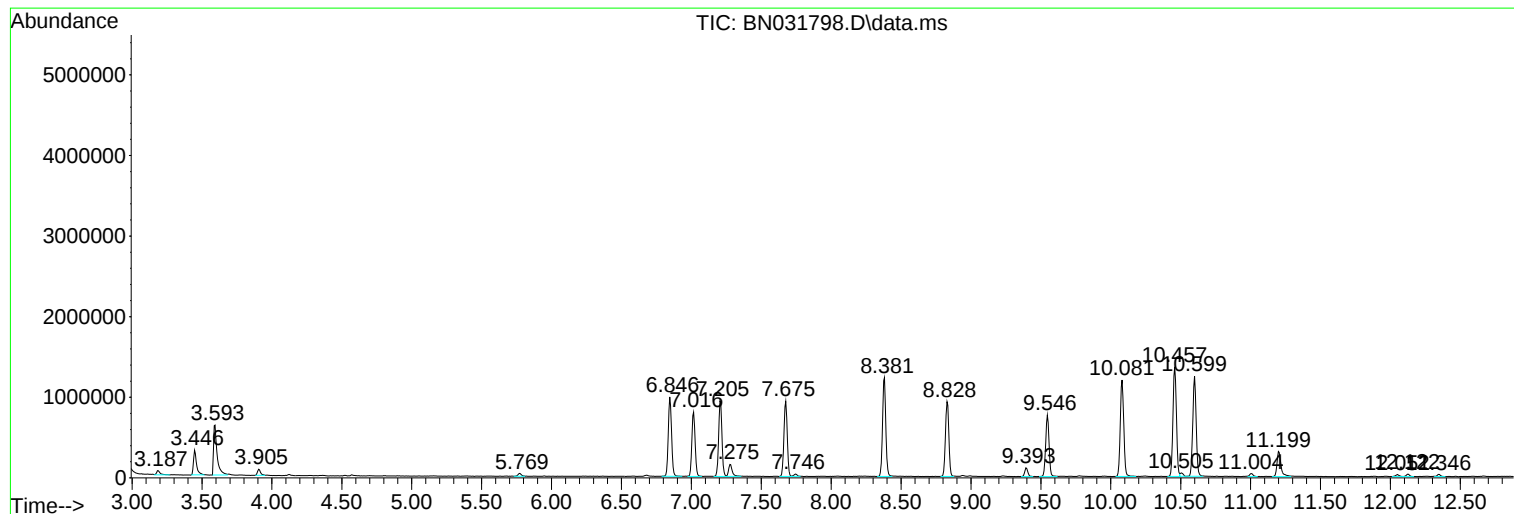
Sum of corrected areas: 108012547

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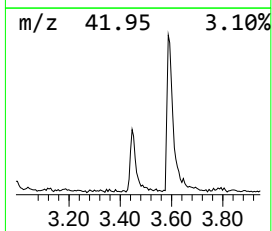
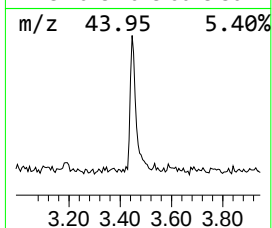
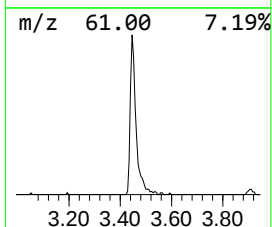
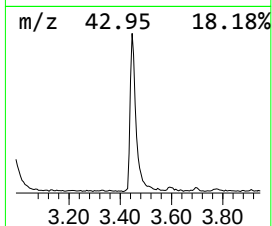
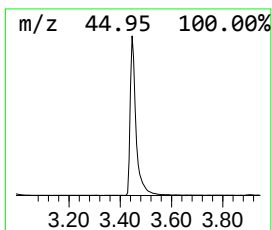
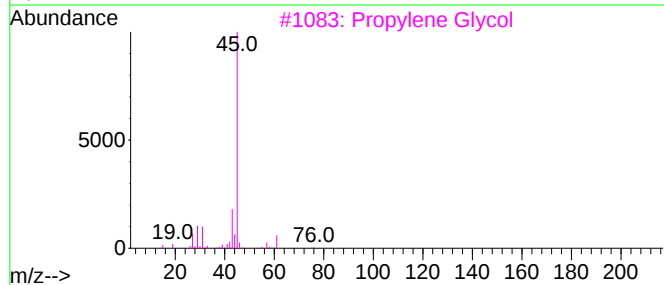
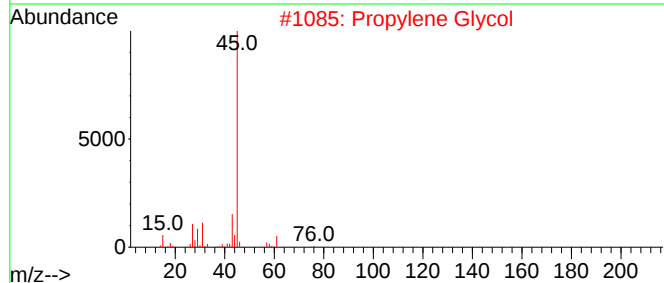
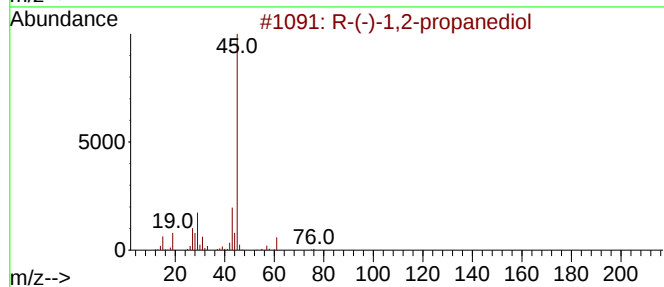
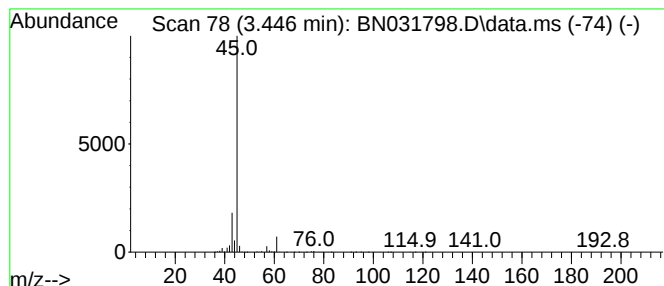
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN052924.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 R-(-)-1,2-propanediol Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.446	6.18 ng/ul	467515	1,4-Dichlorobenzene-d4	7.675

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	R-(-)-1,2-propanediol			76	C3H8O2	004254-14-2	72
2	Propylene Glycol			76	C3H8O2	000057-55-6	64
3	Propylene Glycol			76	C3H8O2	000057-55-6	49
4	Isopropyl Alcohol			60	C3H8O	000067-63-0	9
5	Isopropyl Alcohol			60	C3H8O	000067-63-0	9



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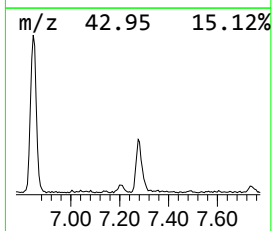
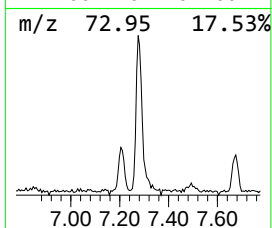
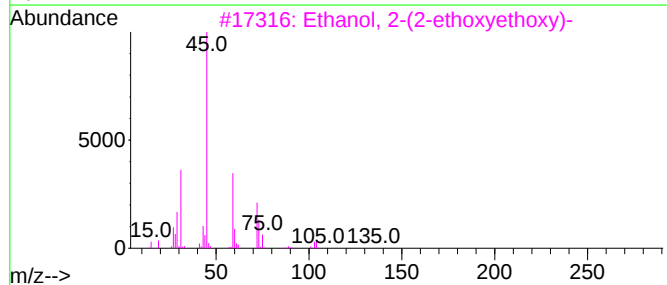
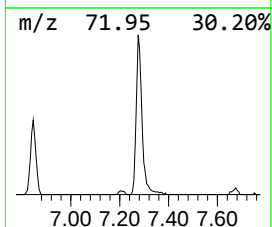
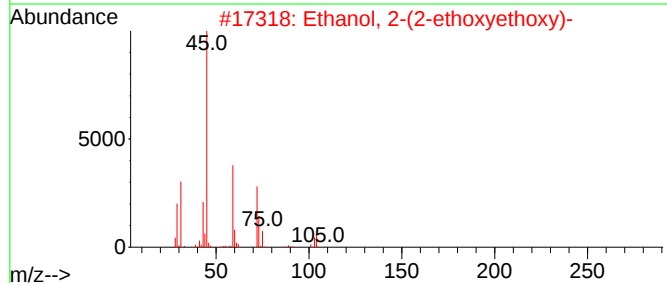
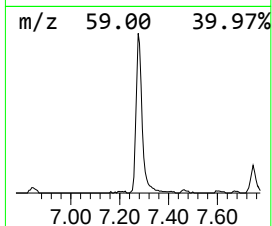
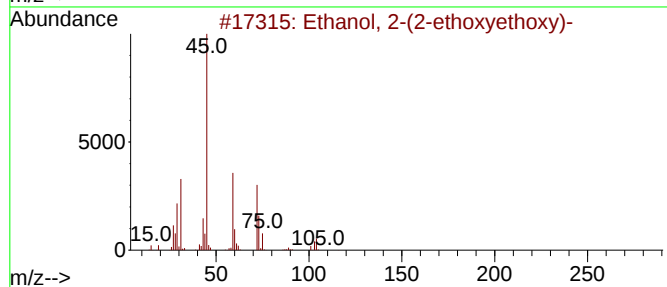
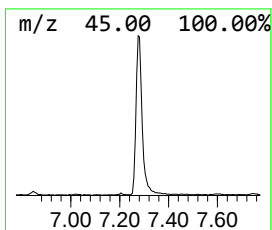
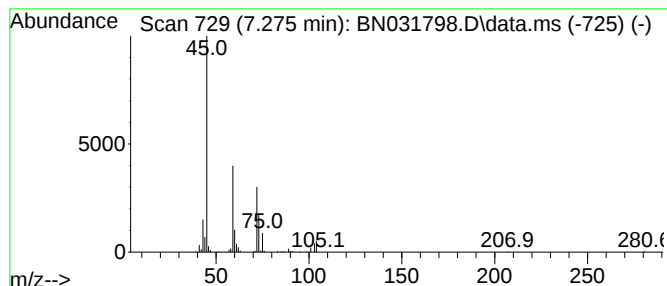
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN052924.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Ethanol, 2-(2-ethoxyethoxy)- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.275	3.48 ng/ul	263524	1,4-Dichlorobenzene-d4	7.675

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	90
4			Ethanol, 2-(2-ethoxyethoxy)-	134	C6H14O3	000111-90-0	90
5			Diethyl carbitol	162	C8H18O3	000112-36-7	72



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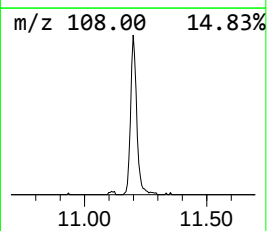
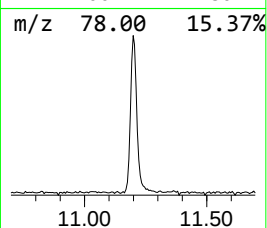
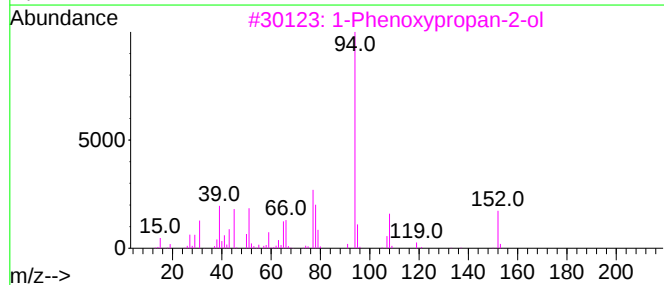
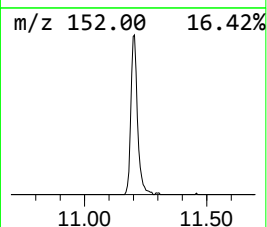
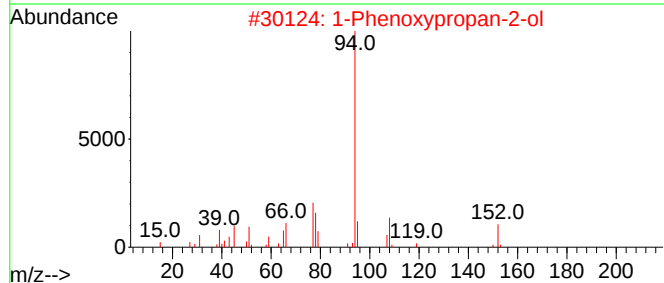
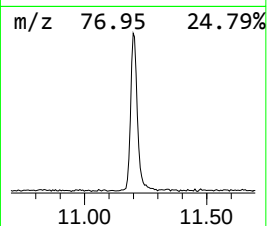
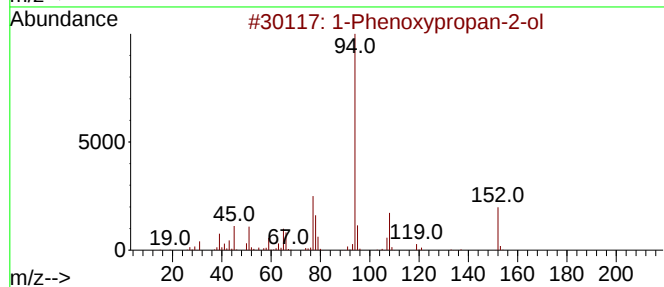
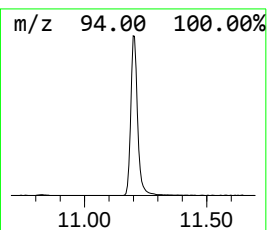
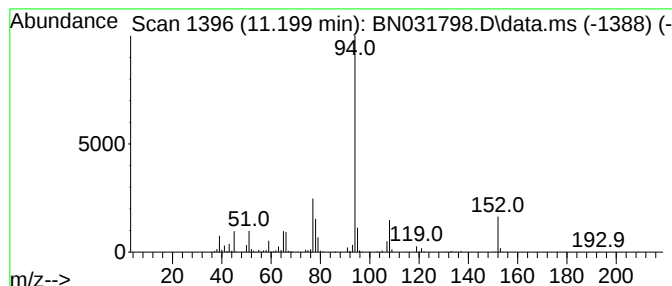
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.199	5.32 ng/ul	623508	Naphthalene-d8	10.457

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-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	95
4			1-Propanol, 3-phenoxy-	152	C9H12O2	006180-61-6	91
5			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN060524\
Data File : BN031798.D
Acq On : 05 Jun 2024 13:48
Operator : MA/JU
Sample : P2623-04
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
BHBQ9

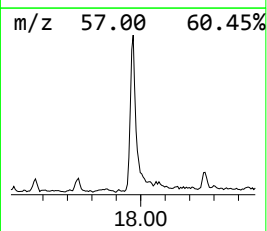
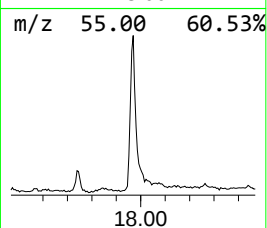
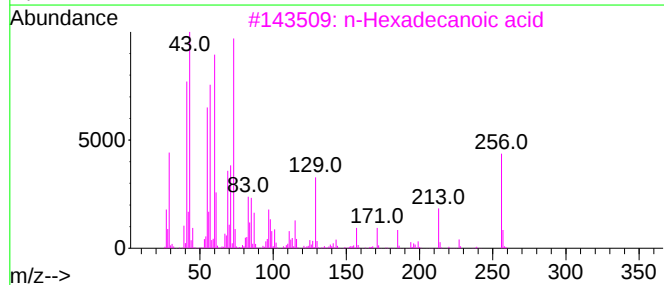
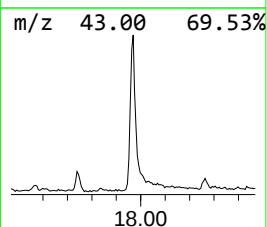
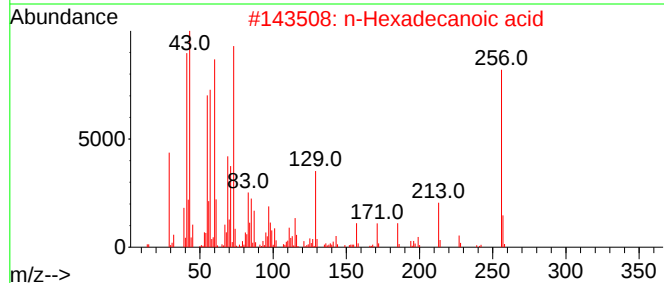
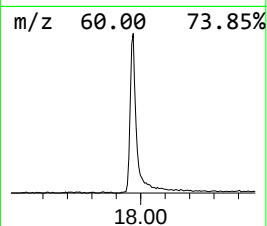
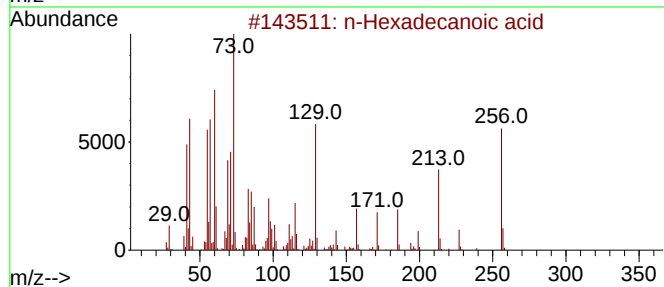
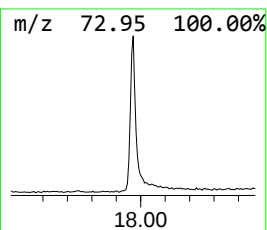
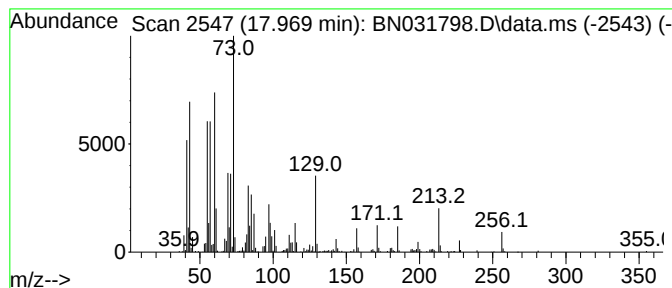
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 n-Hexadecanoic acid Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.969	5.95 ng/ul	1268910	Phenanthrene-d10	17.069

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	96
5			Tridecanoic acid	214	C13H26O2	000638-53-9	89



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN060524\
Data File : BN031798.D
Acq On : 05 Jun 2024 13:48
Operator : MA/JU
Sample : P2623-04
Misc :
ALS Vial : 4 Sample Multiplier: 1

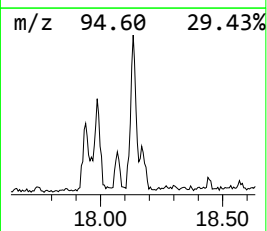
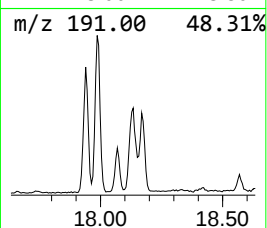
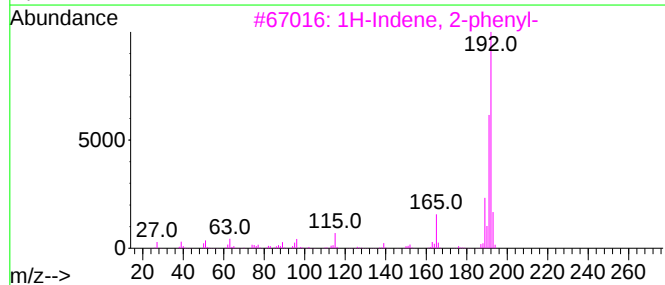
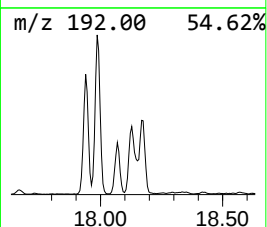
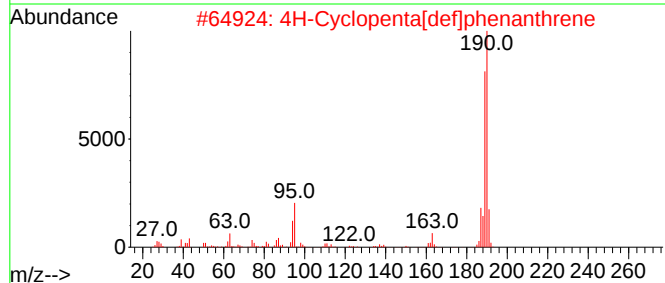
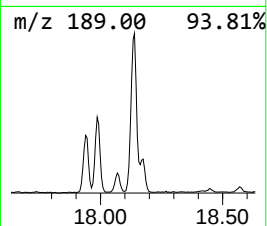
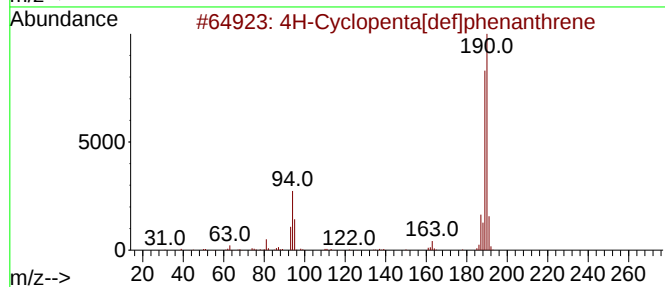
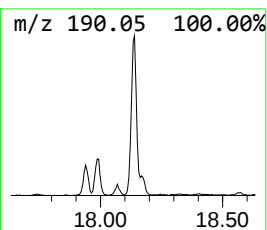
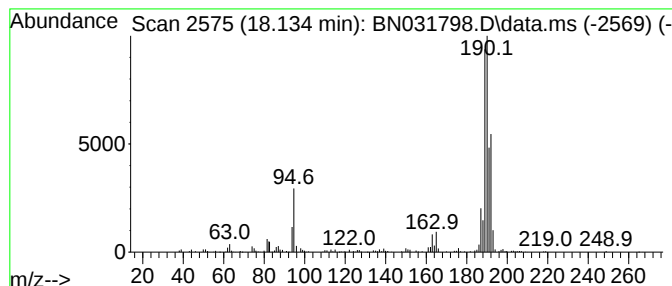
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN052924.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 5 4H-Cyclopenta[def]phenanthrene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
18.134	2.23 ng/ul	474867	Phenanthrene-d10	17.069	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	93
2	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	55
3	1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	35
4	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	25
5	Cyclohexanone, 2-(2-furanylmethy...	190	C12H14O2	056053-07-7	22



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN060524\
Data File : BN031798.D
Acq On : 05 Jun 2024 13:48
Operator : MA/JU
Sample : P2623-04
Misc :
ALS Vial : 4 Sample Multiplier: 1

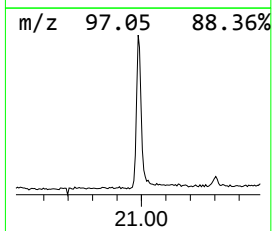
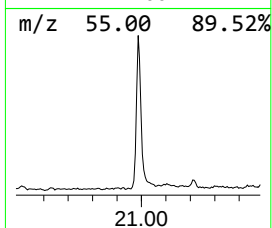
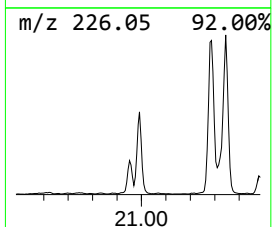
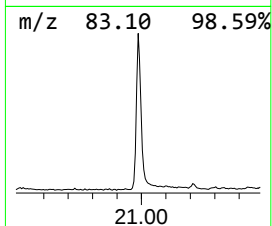
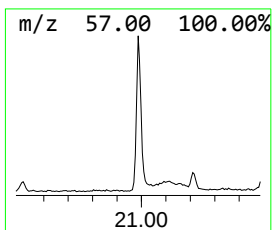
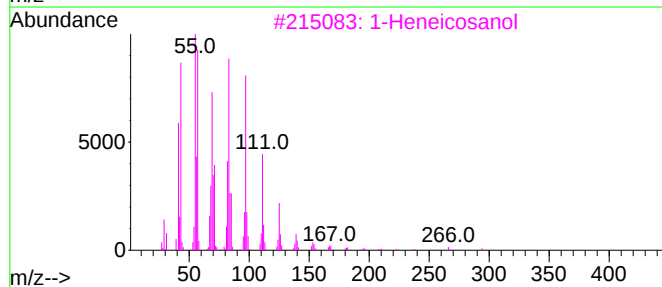
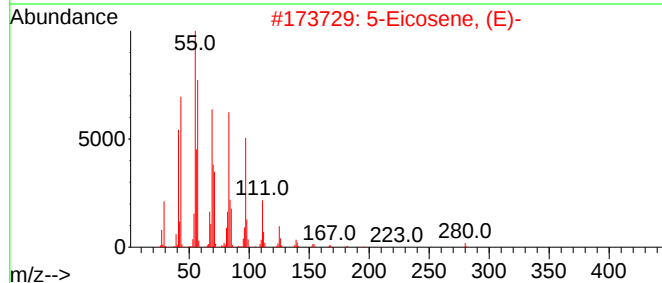
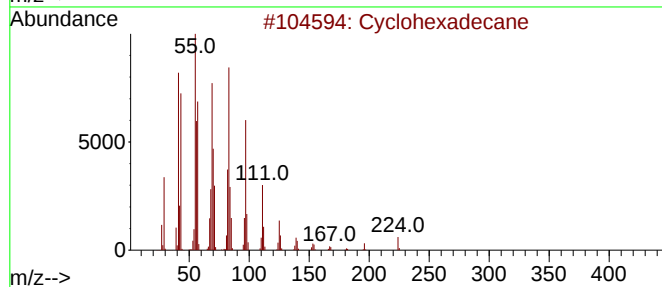
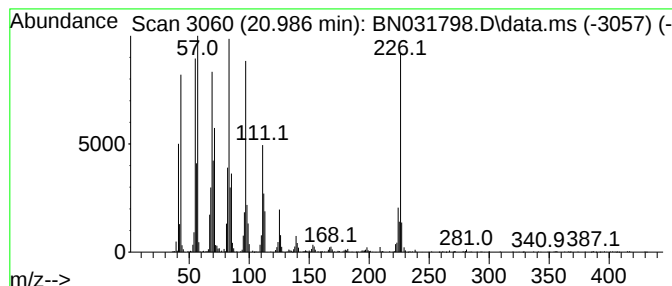
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 (DEL) Alkane: Cyclic20.986 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.	
20.986	3.61 ng/ul	1409680	Chrysene-d12	21.304	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexadecane	224	C16H32	000295-65-8	98
2	5-Eicosene, (E)-	280	C20H40	074685-30-6	97
3	1-Heneicosanol	312	C21H44O	015594-90-8	93
4	1-Heneicosyl formate	340	C22H44O2	077899-03-7	89
5	Trifluoroacetic acid, pentadecyl...	324	C17H31F3O2	959010-23-2	89



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN060524\
Data File : BN031798.D
Acq On : 05 Jun 2024 13:48
Operator : MA/JU
Sample : P2623-04
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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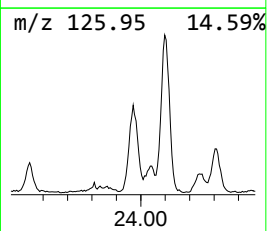
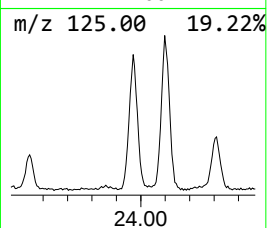
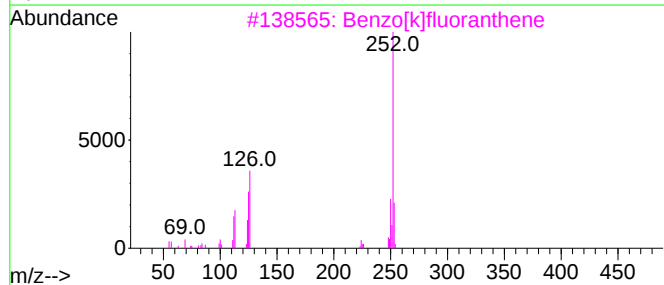
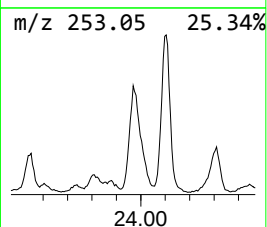
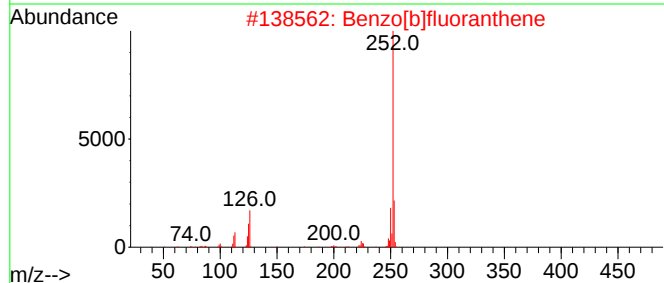
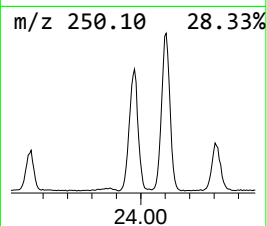
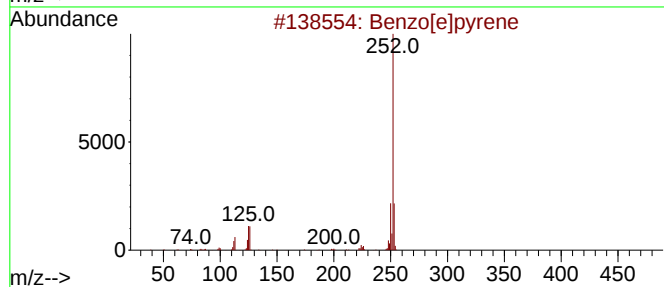
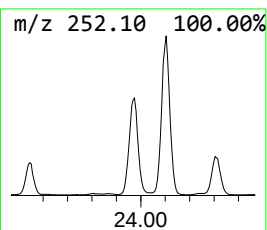
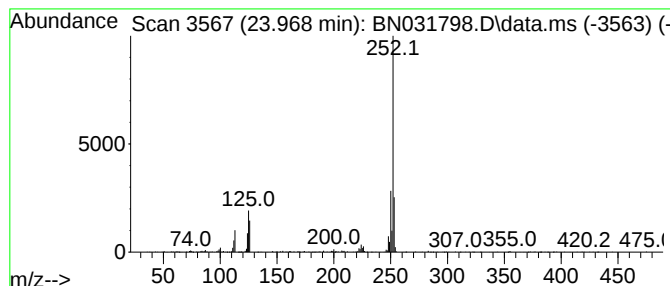
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Benzo[e]pyrene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.968	2.21 ng/ul	673904	Perylene-d12	24.245

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo[e]pyrene	252	C20H12	000192-97-2	98
2			Benzo[b]fluoranthene	252	C20H12	000205-99-2	97
3			Benzo[k]fluoranthene	252	C20H12	000207-08-9	95
4			Benzo[b]fluoranthene	252	C20H12	000205-99-2	94
5			Benzo[a]pyrene	252	C20H12	000050-32-8	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN060524\
Data File : BN031798.D
Acq On : 05 Jun 2024 13:48
Operator : MA/JU
Sample : P2623-04
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
BHBQ9

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
R-(-)-1,2-propa...	3.446	6.2	ng/ul	467515	1	7.675	1514220	20.0
Ethanol, 2-(2-e...	7.275	3.5	ng/ul	263524	1	7.675	1514220	20.0
1-Phenoxypropan...	11.199	5.3	ng/ul	623508	2	10.457	2343620	20.0
n-Hexadecanoic ...	17.969	6.0	ng/ul	1268910	4	17.069	4262680	20.0
4H-Cyclopenta[d...	18.134	2.2	ng/ul	474867	4	17.069	4262680	20.0
(DEL) Alkane: C...	20.986	3.6	ng/ul	1409680	5	21.304	7804580	20.0
Benzo[e]pyrene	23.968	2.2	ng/ul	673904	6	24.245	6112250	20.0