

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062024\
 Data File : BN032109.D
 Acq On : 20 Jun 2024 23:17
 Operator : MA/JU
 Sample : P2710-08
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A4BL7

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: BN032109.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.058	9	12	23	rVB7	11064	26792	0.35%	0.025%
2	3.164	26	30	40	rVB	57284	82165	1.07%	0.077%
3	3.434	72	76	84	rBV	34306	52848	0.69%	0.050%
4	3.575	98	100	110	rBV	86373	127856	1.67%	0.120%
5	3.669	112	116	124	rVB	69315	99322	1.29%	0.093%
6	3.881	148	152	156	rVB	17986	25108	0.33%	0.024%
7	4.375	231	236	240	rBV2	22957	33051	0.43%	0.031%
8	4.422	240	244	251	rVB	47865	71732	0.94%	0.067%
9	4.769	298	303	311	rBV	133336	205507	2.68%	0.193%
10	6.322	562	567	574	rVB3	15289	24355	0.32%	0.023%
11	6.652	614	623	635	rBV2	24055	55825	0.73%	0.052%
12	6.816	644	651	668	rVB	1170954	1918341	25.01%	1.798%
13	6.981	672	679	690	rBV	980665	1514170	19.74%	1.419%
14	7.169	703	711	719	rBV	1228177	2005536	26.14%	1.879%
15	7.246	719	724	736	rVB	135976	236681	3.09%	0.222%
16	7.634	782	790	797	rBV	1216395	1961492	25.57%	1.838%
17	7.699	797	801	810	rVB4	14224	30973	0.40%	0.029%
18	8.346	902	911	926	rBV	1475132	2457283	32.03%	2.303%
19	8.463	926	931	938	rVB	64613	111798	1.46%	0.105%
20	8.793	979	987	996	rBV	1189400	1962537	25.58%	1.839%
21	9.357	1078	1083	1096	rVB	25626	45433	0.59%	0.043%
22	9.510	1101	1109	1120	rBV	999074	1699747	22.16%	1.593%
23	9.746	1141	1149	1154	rVB5	14723	35328	0.46%	0.033%
24	10.046	1191	1200	1218	rBV	1569675	2836758	36.98%	2.658%
25	10.416	1255	1263	1277	rBV	1897148	3232580	42.14%	3.029%
26	10.563	1281	1288	1307	rBV	788255	1479659	19.29%	1.387%
27	10.963	1348	1356	1366	rBV9	15826	44151	0.58%	0.041%
28	11.163	1384	1390	1396	rBV	79522	157216	2.05%	0.147%
29	11.210	1396	1398	1411	rVB3	23781	48767	0.64%	0.046%
30	12.010	1527	1534	1540	rVV	31855	57888	0.75%	0.054%
31	12.634	1634	1640	1646	rBV8	14977	38289	0.50%	0.036%
32	13.175	1724	1732	1741	rBV	802914	1299868	16.94%	1.218%
33	13.592	1798	1803	1810	rVB7	16219	32332	0.42%	0.030%
34	13.698	1814	1821	1834	rVV	3169578	4347013	56.67%	4.074%
35	13.975	1856	1868	1885	rBV2	3366896	5414512	70.58%	5.074%
36	14.116	1885	1892	1896	rBV4	27683	49782	0.65%	0.047%

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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	14.281	1910	1920	1927	rBV	3103052	4649519	60.61%	4.357%
38	14.345	1927	1931	1937	rVB	53983	82561	1.08%	0.077%
39	14.404	1937	1941	1947	rVB2	22413	34728	0.45%	0.033%
40	14.504	1950	1958	1970	rBV	1181214	2064567	26.91%	1.935%
41	14.657	1979	1984	1990	rBV2	86847	163349	2.13%	0.153%
42	14.710	1990	1993	1998	rVB4	36007	54351	0.71%	0.051%
43	15.081	2050	2056	2059	rBV2	23733	40546	0.53%	0.038%
44	15.281	2082	2090	2095	rVV	4426947	6574950	85.71%	6.162%
45	15.334	2095	2099	2106	rVV	62194	104492	1.36%	0.098%
46	15.410	2106	2112	2123	rVV	1206937	1726031	22.50%	1.618%
47	15.539	2125	2134	2140	rVV2	90003	193290	2.52%	0.181%
48	15.628	2144	2149	2159	rVV7	24577	65397	0.85%	0.061%
49	15.775	2171	2174	2181	rVB4	24058	43218	0.56%	0.041%
50	15.839	2181	2185	2193	rBV4	16609	37058	0.48%	0.035%
51	16.004	2208	2213	2220	rVB9	31161	70608	0.92%	0.066%
52	16.181	2239	2243	2252	rVB5	75607	124764	1.63%	0.117%
53	16.263	2253	2257	2261	rBV2	55360	70781	0.92%	0.066%
54	16.334	2267	2269	2275	rVB6	18898	24600	0.32%	0.023%
55	16.439	2282	2287	2293	rBV3	66492	116955	1.52%	0.110%
56	16.563	2301	2308	2313	rBV7	43081	98575	1.28%	0.092%
57	16.686	2325	2329	2337	rVB2	38803	69766	0.91%	0.065%
58	16.757	2337	2341	2343	rBV4	23172	35451	0.46%	0.033%
59	16.792	2344	2347	2353	rVV4	39174	75625	0.99%	0.071%
60	16.845	2353	2356	2362	rVB	54468	84689	1.10%	0.079%
61	16.945	2364	2373	2378	rBV6	45049	111265	1.45%	0.104%
62	17.028	2381	2387	2392	rBV	3614771	5486913	71.52%	5.142%
63	17.069	2392	2394	2399	rVV	933420	1229265	16.02%	1.152%
64	17.128	2399	2404	2414	rVV2	4466752	6907018	90.04%	6.473%
65	17.210	2414	2418	2425	rVB2	44825	99761	1.30%	0.093%
66	17.439	2446	2457	2463	rBV2	104121	254767	3.32%	0.239%
67	17.616	2483	2487	2494	rVB5	38702	89488	1.17%	0.084%
68	17.704	2498	2502	2506	rBV	115231	142893	1.86%	0.134%
69	17.745	2506	2509	2515	rVB4	30578	47310	0.62%	0.044%
70	17.810	2515	2520	2527	rBV7	46410	111181	1.45%	0.104%
71	17.933	2528	2541	2550	rBV2	1805874	3224932	42.04%	3.022%
72	18.098	2564	2569	2573	rBV2	218678	352621	4.60%	0.330%
73	18.133	2573	2575	2579	rVB	64728	77003	1.00%	0.072%
74	18.263	2594	2597	2601	rVB3	43045	52699	0.69%	0.049%
75	18.410	2618	2622	2626	rBV	118069	168069	2.19%	0.158%
76	18.457	2626	2630	2635	rVB	138152	185256	2.41%	0.174%

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If leading or trailing edge < 100 prefer < Baseline drop else tangent >

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 Title : SVOA CALIBRATION

77	18.657	2661	2664	2669	rVB4	52322	66999	0.87%	0.063%
78	18.828	2690	2693	2696	rBV2	55693	77579	1.01%	0.073%
79	18.857	2696	2698	2700	rVV2	72642	90180	1.18%	0.085%
80	18.886	2700	2703	2707	rVB3	185394	225955	2.95%	0.212%
81	18.986	2714	2720	2724	rBV3	94484	212556	2.77%	0.199%
82	19.092	2730	2738	2745	rBV	2556445	3714773	48.42%	3.481%
83	19.222	2756	2760	2769	rBV2	751962	1062003	13.84%	0.995%
84	19.427	2790	2795	2799	rBV	4703547	7671344	100.00%	7.189%
85	19.527	2809	2812	2817	rBV2	88010	125583	1.64%	0.118%
86	19.639	2828	2831	2836	rVB	104754	133091	1.73%	0.125%
87	19.786	2851	2856	2862	rBV4	116322	270344	3.52%	0.253%
88	19.957	2881	2885	2887	rBV2	270192	339478	4.43%	0.318%
89	20.057	2898	2902	2906	rBV	194891	259702	3.39%	0.243%
90	20.474	2970	2973	2976	rVB	215545	238531	3.11%	0.224%
91	20.951	3050	3054	3061	rVB2	1286569	1853592	24.16%	1.737%
92	21.169	3088	3091	3096	rVB	426944	499393	6.51%	0.468%
93	21.263	3099	3107	3111	rVV2	3211171	6084836	79.32%	5.702%
94	21.304	3111	3114	3125	rVB	950149	1355219	17.67%	1.270%
95	21.998	3228	3232	3235	rBV2	487822	788924	10.28%	0.739%
96	23.251	3439	3445	3451	rBV	570260	1491871	19.45%	1.398%
97	23.333	3453	3459	3466	rVB2	1331529	2939017	38.31%	2.754%
98	23.968	3561	3567	3574	rBV2	1294342	2876470	37.50%	2.696%
99	24.174	3594	3602	3612	rVB	1732914	4217739	54.98%	3.953%
100	25.174	3766	3772	3783	rVB2	429014	1142231	14.89%	1.070%

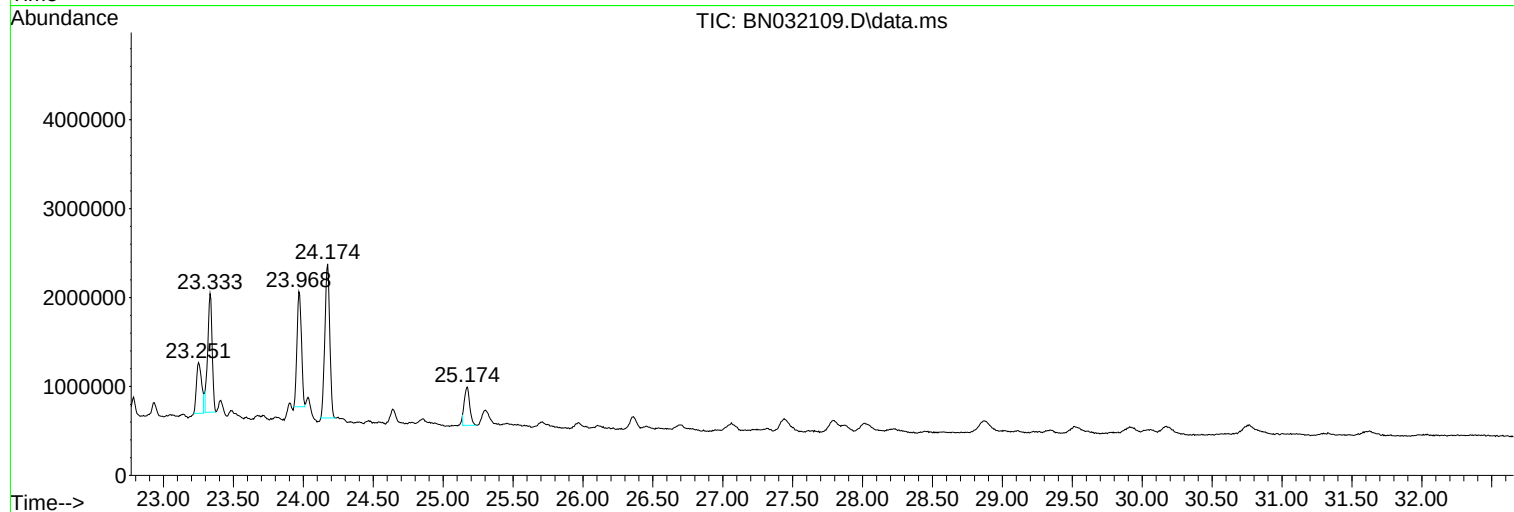
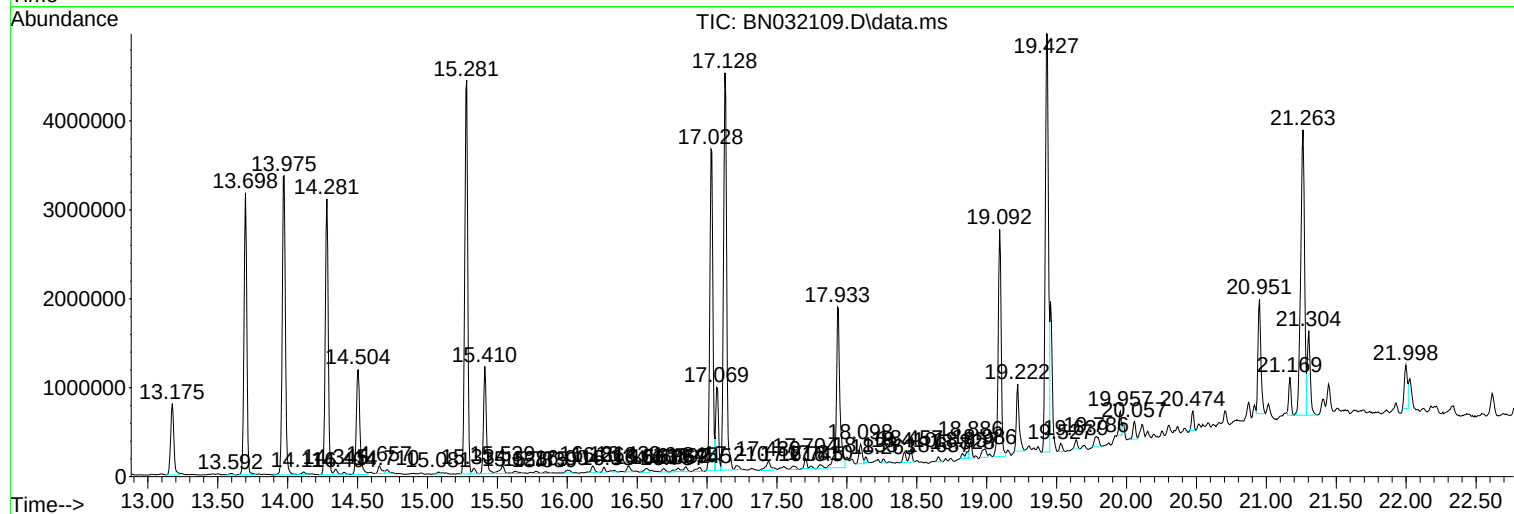
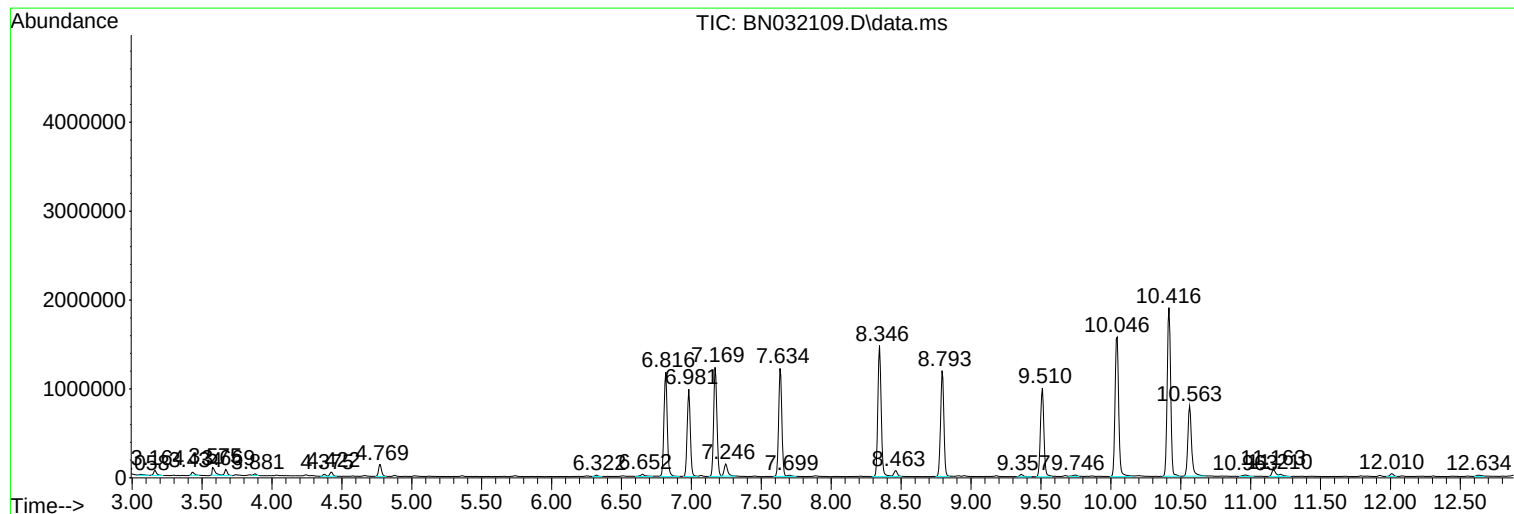
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Instrument :
BNA_N
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A4BL7

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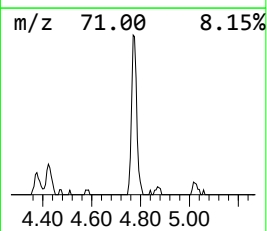
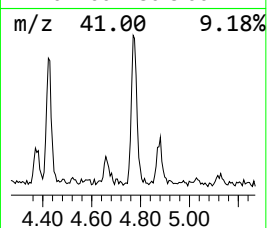
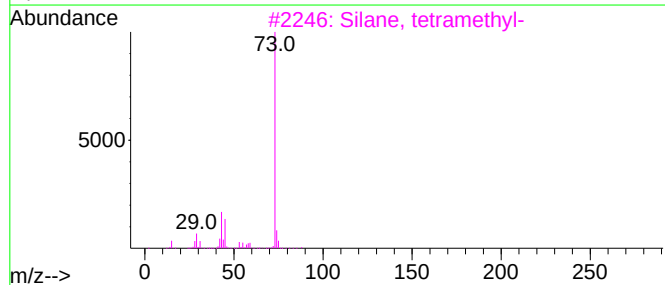
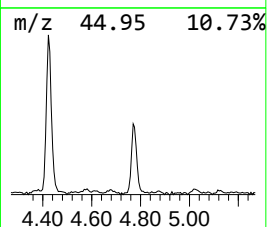
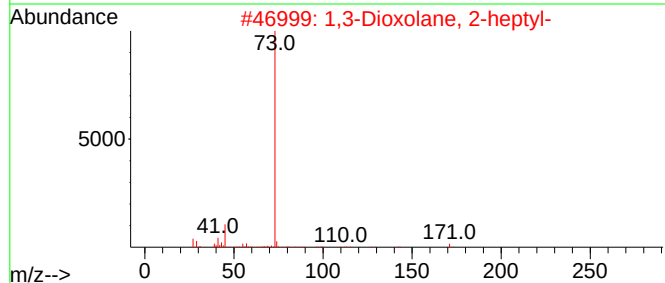
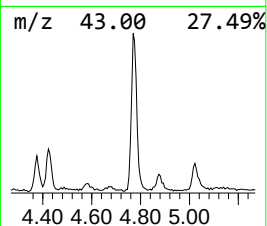
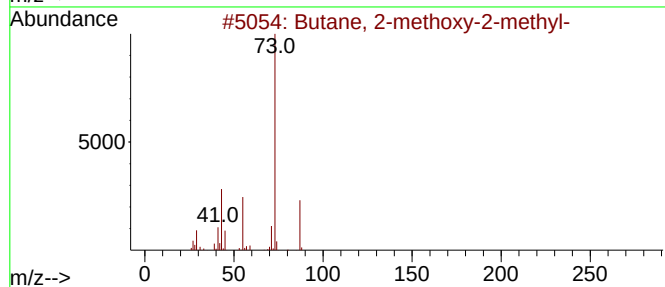
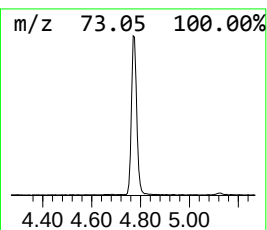
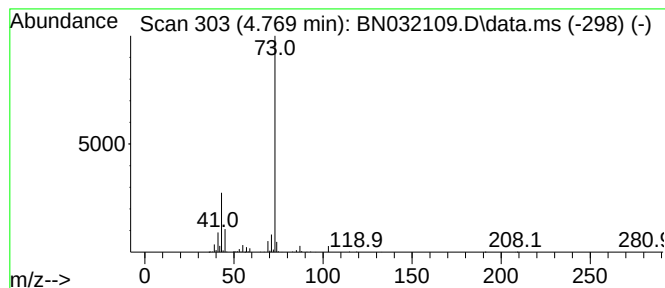
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN052924.MA.M
Quant Title : SVOA CALIBRATION

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	2.10 ng/ul	205507	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			1,3-Dioxolane, 2-heptyl-	172	C10H20O2	004359-57-3	38
3			Silane, tetramethyl-	88	C4H12Si	000075-76-3	37
4			1-Propene-1-thiol	74	C3H6S	000925-89-3	9
5			N-Ethylformamide	73	C3H7NO	000627-45-2	9



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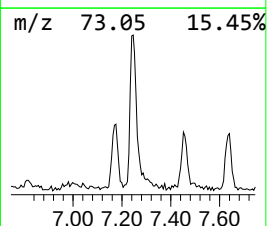
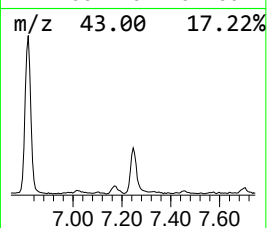
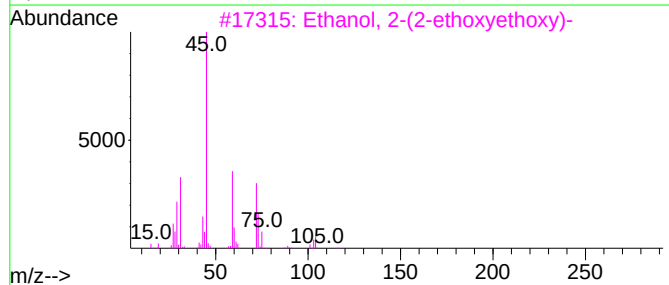
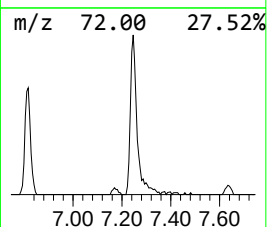
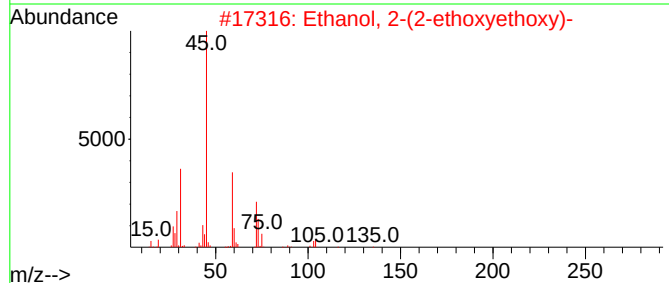
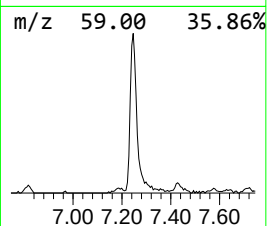
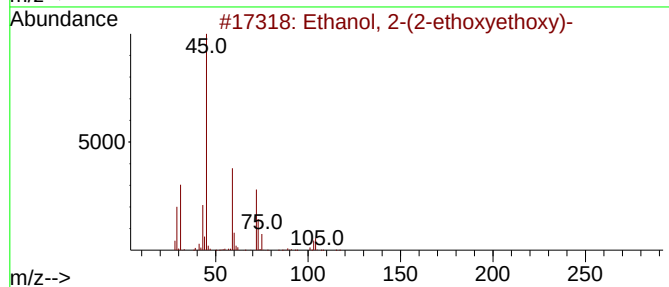
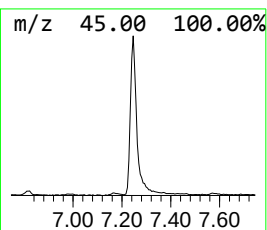
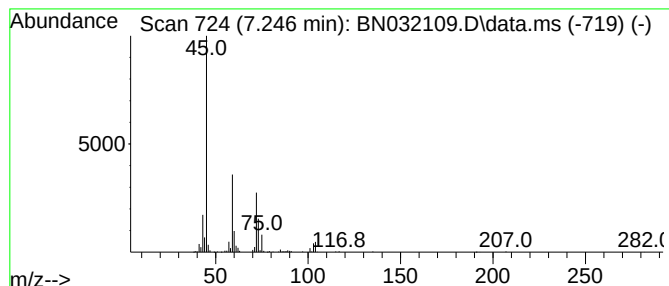
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN052924.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Ethanol, 2-(2-ethoxyethoxy)- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.246	2.41 ng/ul	236681	1,4-Dichlorobenzene-d4	7.634

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	83
4			Diethyl carbitol	162	C8H18O3	000112-36-7	72
5			Ethanol, 2-(2-ethoxyethoxy)-	134	C6H14O3	000111-90-0	50



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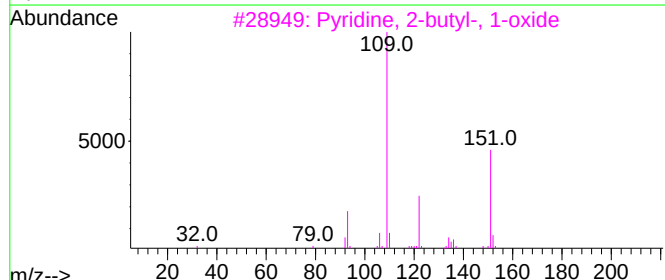
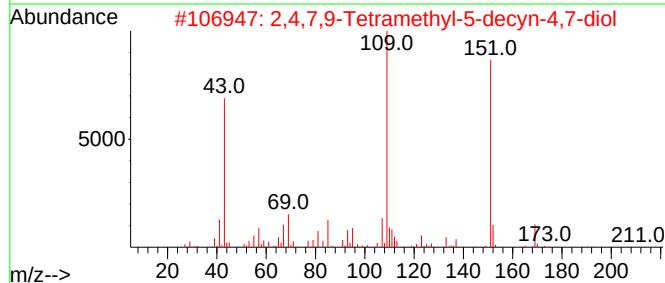
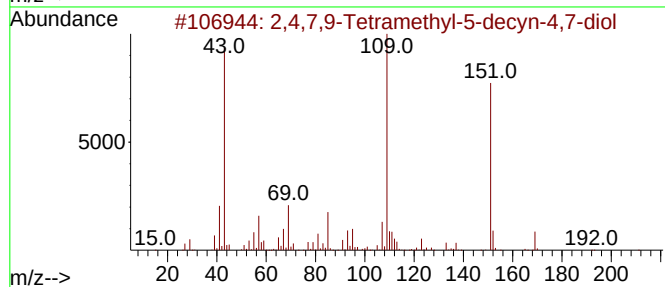
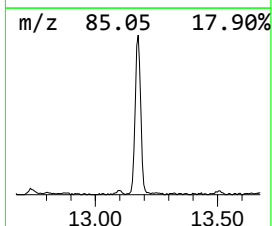
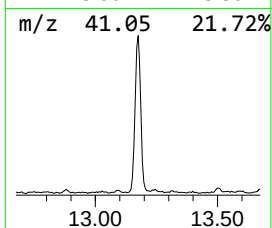
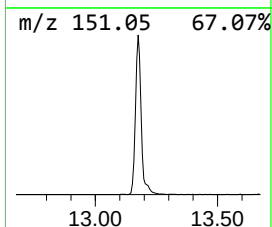
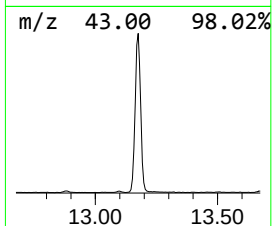
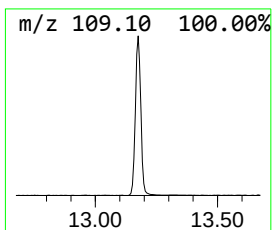
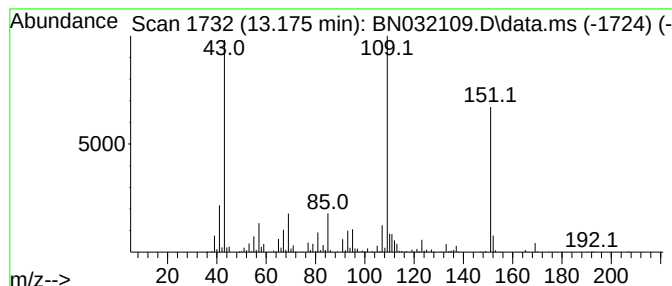
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN052924.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 2,4,7,9-Tetramethyl-5-decyn... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.175	5.59 ng/ul	1299870	Acenaphthene-d10	14.281

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,4,7,9-Tetramethyl-5-decyn-4,7-...	226	C14H26O2	000126-86-3	91		
2	2,4,7,9-Tetramethyl-5-decyn-4,7-...	226	C14H26O2	000126-86-3	91		
3	Pyridine, 2-butyl-, 1-oxide	151	C9H13NO	031396-32-4	59		
4	Metacetamol	151	C8H9NO2	000621-42-1	53		
5	2-Propenal, 3-(2,2,6-trimethyl-7...	194	C12H18O2	055759-91-6	50		



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062024\
Data File : BN032109.D
Acq On : 20 Jun 2024 23:17
Operator : MA/JU
Sample : P2710-08
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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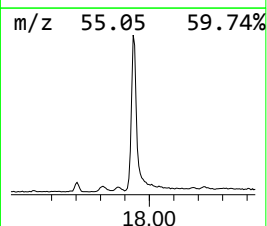
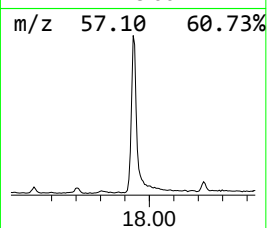
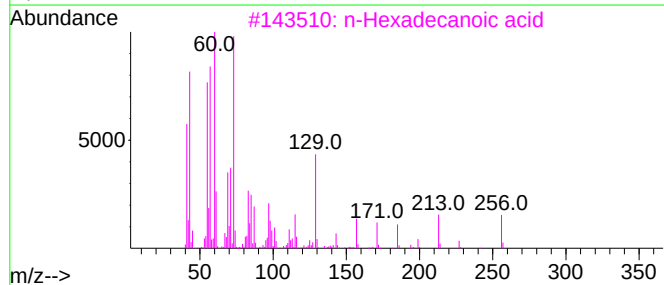
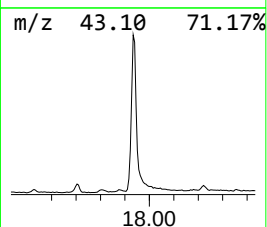
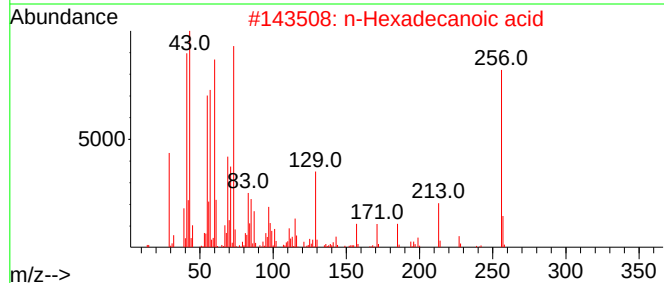
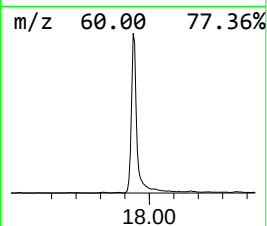
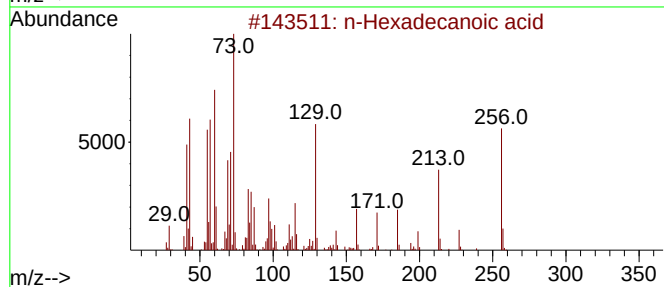
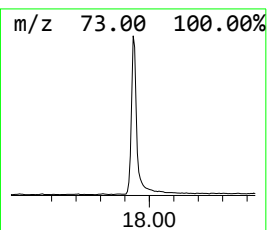
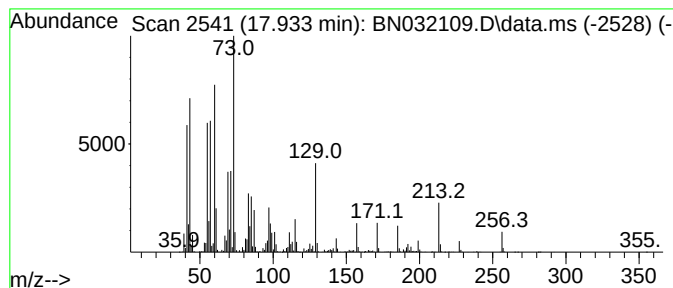
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 n-Hexadecanoic acid Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.933	11.76 ng/ul	3224930	Phenanthrene-d10	17.028

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	99
5			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062024\
Data File : BN032109.D
Acq On : 20 Jun 2024 23:17
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Sample : P2710-08
Misc :
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Instrument :
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ClientSampleId :
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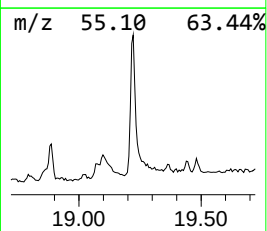
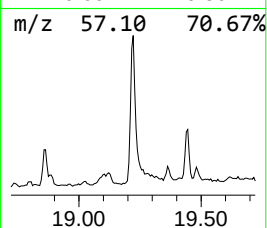
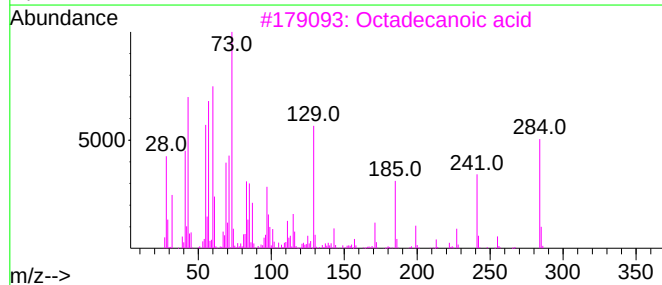
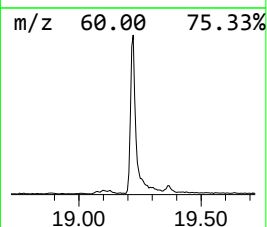
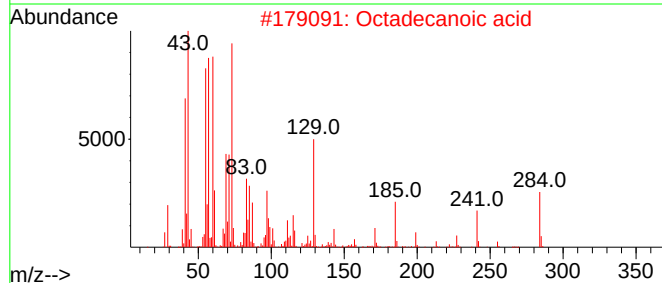
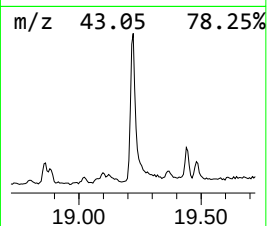
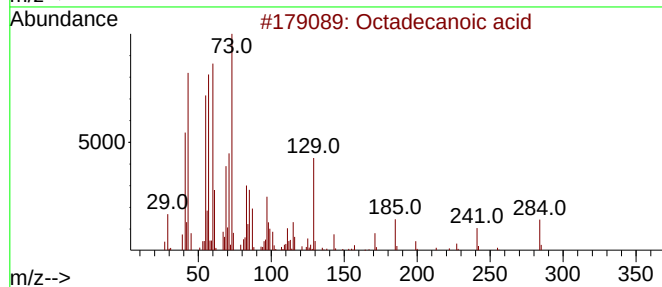
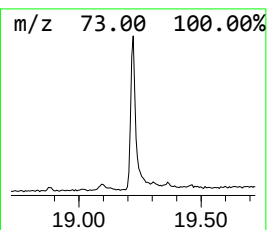
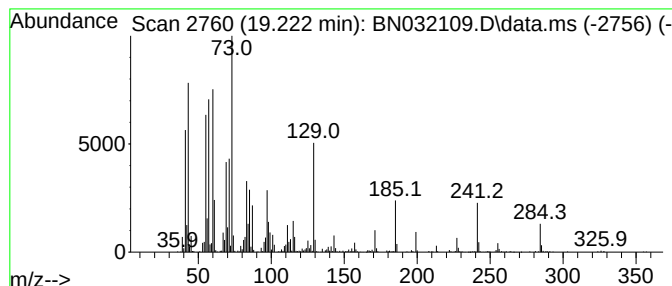
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Octadecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.222	3.49 ng/ul	1062000	Chrysene-d12	21.263

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\BN062024\
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Operator : MA/JU
Sample : P2710-08
Misc :
ALS Vial : 17 Sample Multiplier: 1

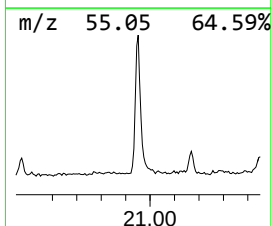
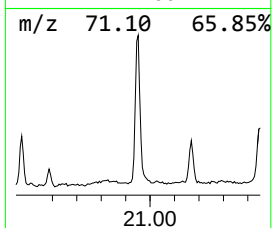
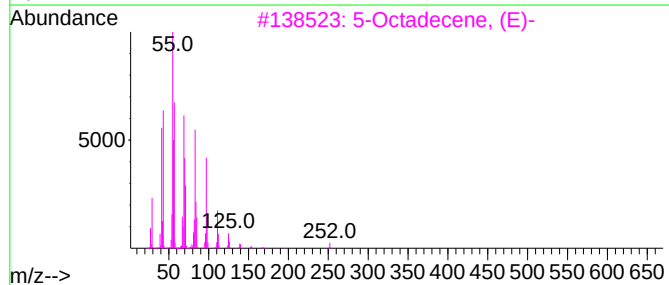
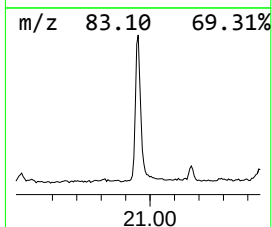
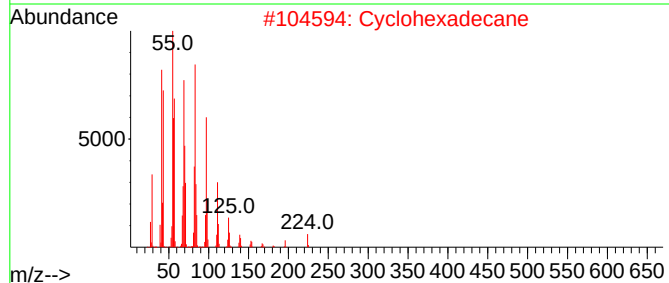
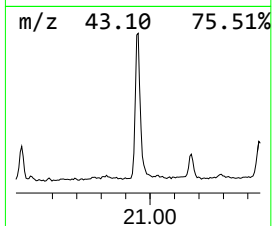
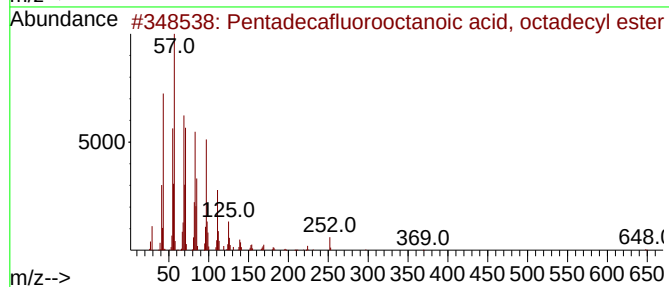
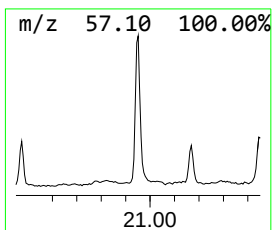
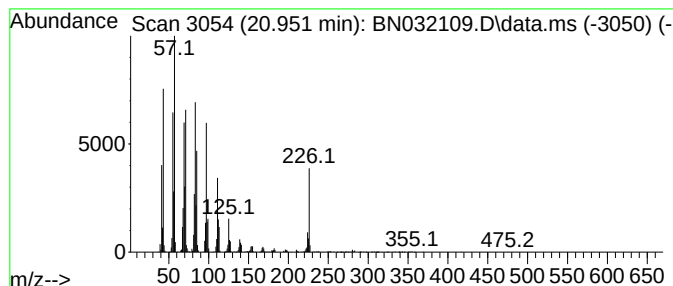
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN052924.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 6 Pentadecafluorooctanoic aci... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
20.951	6.09 ng/ul	1853590	Chrysene-d12	21.263	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentadecafluorooctanoic acid, oc...	666	C26H37F15O2	1000406-04-8	93
2	Cyclohexadecane	224	C16H32	000295-65-8	91
3	5-Octadecene, (E)-	252	C18H36	007206-21-5	91
4	1-Heneicosanol	312	C21H44O	015594-90-8	90
5	Pentadecafluorooctanoic acid, he...	638	C24H33F15O2	1000406-04-6	89



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062024\
Data File : BN032109.D
Acq On : 20 Jun 2024 23:17
Operator : MA/JU
Sample : P2710-08
Misc :
ALS Vial : 17 Sample Multiplier: 1

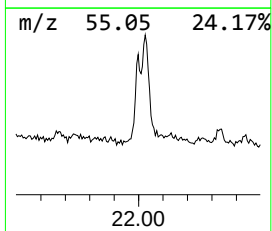
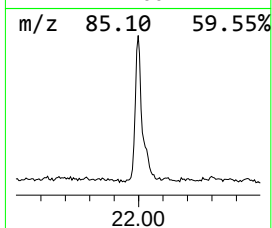
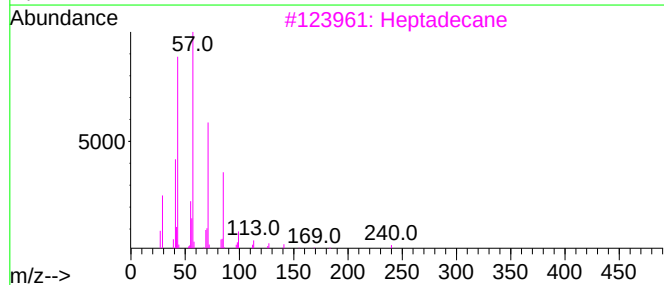
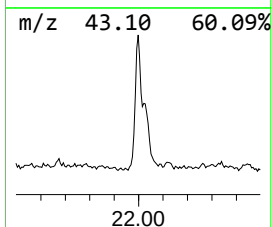
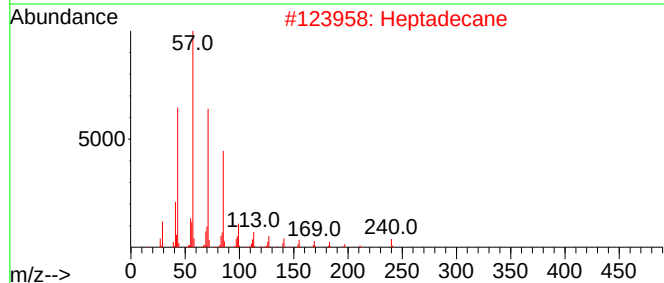
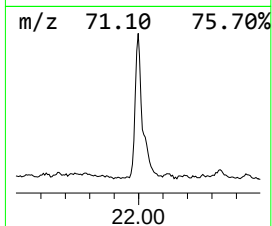
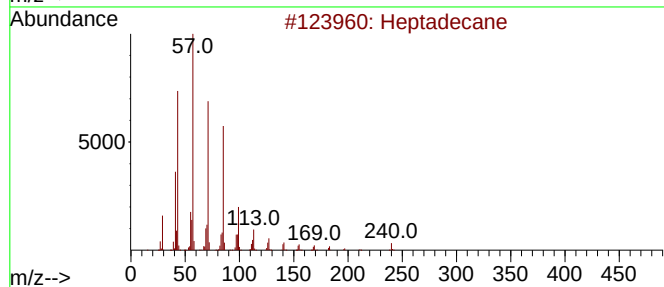
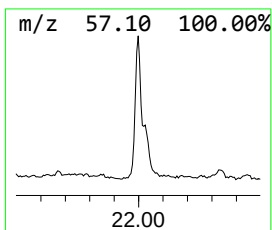
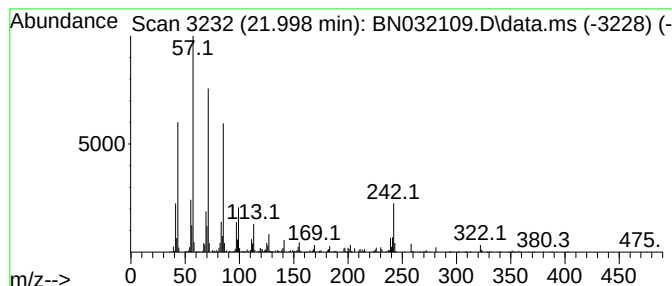
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD		R.T.	
21.998	2.59 ng/ul	788924	Chrysene-d12		21.263	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Heptadecane		240	C17H36	000629-78-7	97
2	Heptadecane		240	C17H36	000629-78-7	95
3	Heptadecane		240	C17H36	000629-78-7	93
4	Pentacosane		352	C25H52	000629-99-2	90
5	Docosane, 11-butyl-		366	C26H54	013475-76-8	89



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062024\
Data File : BN032109.D
Acq On : 20 Jun 2024 23:17
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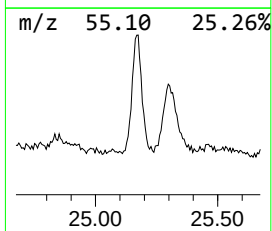
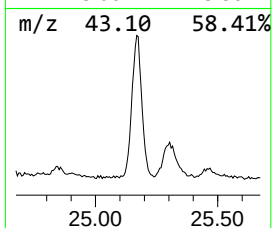
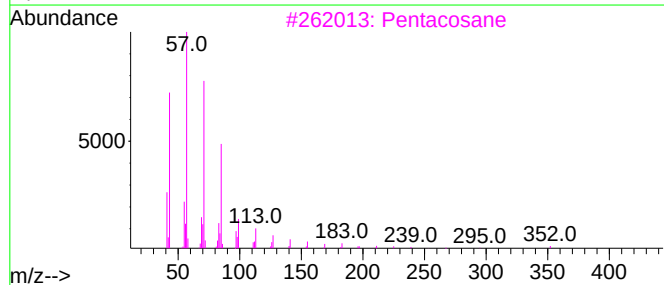
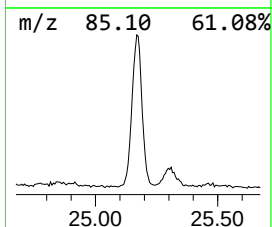
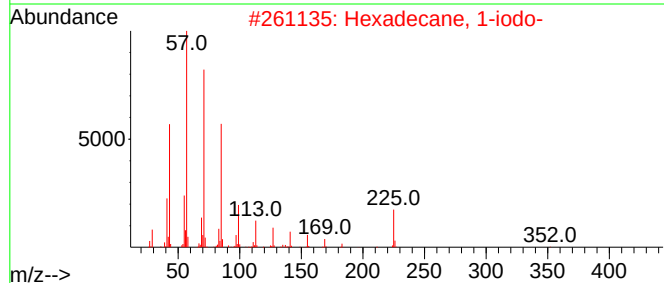
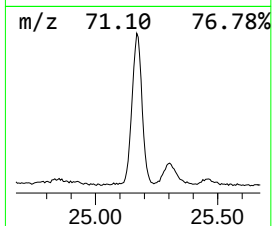
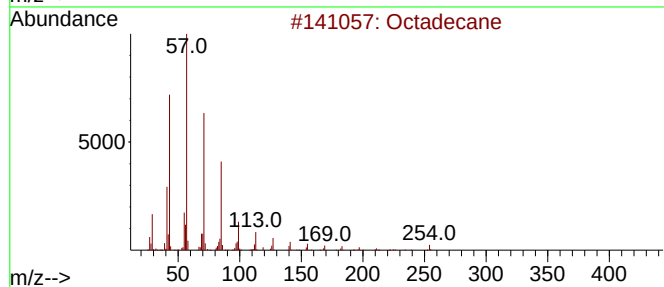
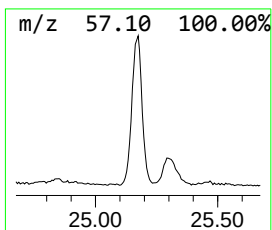
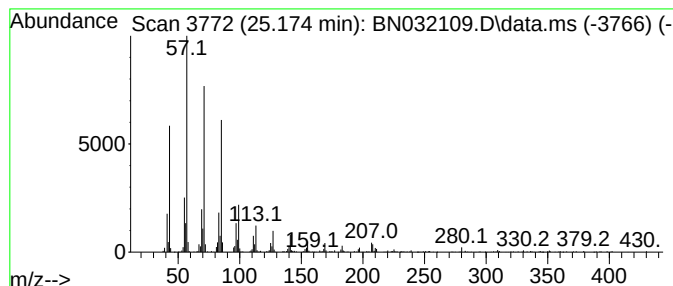
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.174	5.42 ng/ul	1142230	Perylene-d12	24.174

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecane	254	C18H38	000593-45-3	96
2			Hexadecane, 1-iodo-	352	C16H33I	000544-77-4	96
3			Pentacosane	352	C25H52	000629-99-2	94
4			Heneicosane	296	C21H44	000629-94-7	91
5			Octacosane, 2-methyl-	408	C29H60	001560-98-1	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062024\
Data File : BN032109.D
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Operator : MA/JU
Sample : P2710-08
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

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					#	RT	Resp	Conc
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Ethanol, 2-(2-e...	7.246	2.4	ng/ul	236681	1	7.634	1961490	20.0
2,4,7,9-Tetrame...	13.175	5.6	ng/ul	1299870	3	14.281	4649520	20.0
n-Hexadecanoic ...	17.933	11.8	ng/ul	3224930	4	17.028	5486910	20.0
Octadecanoic acid	19.222	3.5	ng/ul	1062000	5	21.263	6084840	20.0
Pentadecafluoro...	20.951	6.1	ng/ul	1853590	5	21.263	6084840	20.0
(DEL) Alkane: S...	21.998	2.6	ng/ul	788924	5	21.263	6084840	20.0
(DEL) Alkane: S...	25.174	5.4	ng/ul	1142230	6	24.174	4217740	20.0