

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062024\
 Data File : BN032112.D
 Acq On : 21 Jun 2024 01:14
 Operator : MA/JU
 Sample : P2710-07
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A4BL4

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: BN032112.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.164	26	30	38	rVB	50555	70328	1.01%	0.063%
2	3.434	73	76	85	rBV	23656	43724	0.62%	0.039%
3	3.581	97	101	112	rBV2	42152	77882	1.11%	0.070%
4	3.670	112	116	124	rVB	80549	112747	1.61%	0.101%
5	4.422	240	244	250	rVB	56388	81200	1.16%	0.073%
6	4.770	298	303	316	rVB	151557	237432	3.39%	0.212%
7	6.652	613	623	633	rBV3	22018	48374	0.69%	0.043%
8	6.816	642	651	664	rVB	1061243	1862117	26.61%	1.665%
9	6.981	672	679	691	rBV	904717	1403862	20.06%	1.255%
10	7.169	703	711	719	rBV	1147966	1857074	26.54%	1.661%
11	7.246	719	724	736	rVV	130664	254267	3.63%	0.227%
12	7.634	783	790	798	rVV	1462385	2352178	33.62%	2.103%
13	7.887	826	833	844	rBV2	31017	71129	1.02%	0.064%
14	8.346	902	911	918	rBV	1414116	2345430	33.52%	2.097%
15	8.399	918	920	926	rVV	42982	73140	1.05%	0.065%
16	8.463	926	931	939	rVB	75016	131024	1.87%	0.117%
17	8.793	980	987	998	rVV	1127521	1827936	26.12%	1.635%
18	9.510	1101	1109	1123	rBV	897853	1580812	22.59%	1.414%
19	10.040	1192	1199	1213	rBV	1490743	2608403	37.28%	2.333%
20	10.416	1254	1263	1269	rBV	2285581	3868366	55.28%	3.459%
21	10.457	1269	1270	1280	rVV2	121388	147518	2.11%	0.132%
22	10.563	1280	1288	1307	rVV	708290	1320867	18.88%	1.181%
23	10.969	1347	1357	1363	rBV	20132	52494	0.75%	0.047%
24	11.163	1384	1390	1395	rBV	76962	151239	2.16%	0.135%
25	11.210	1396	1398	1413	rVB4	25086	55673	0.80%	0.050%
26	11.593	1457	1463	1470	rBV6	26587	49705	0.71%	0.044%
27	12.010	1528	1534	1541	rVB2	28696	52751	0.75%	0.047%
28	13.093	1709	1718	1725	rVB2	17093	43282	0.62%	0.039%
29	13.175	1725	1732	1737	rBV	185423	316686	4.53%	0.283%
30	13.275	1745	1749	1757	rVB	38354	67444	0.96%	0.060%
31	13.440	1771	1777	1785	rBV2	37880	106205	1.52%	0.095%
32	13.698	1809	1821	1835	rBV	2778810	4038845	57.72%	3.612%
33	13.975	1855	1868	1883	rBV	3039887	5006511	71.55%	4.477%
34	14.281	1912	1920	1926	rBV	3691839	5475904	78.26%	4.897%
35	14.346	1926	1931	1938	rVB2	90061	167437	2.39%	0.150%
36	14.504	1951	1958	1970	rBV	932786	1648120	23.55%	1.474%

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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 >
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Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN052924.MA.M
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37	14.657	1979	1984	1987	rBV	78026	132917	1.90%	0.119%
38	15.075	2048	2055	2060	rBV2	27272	62590	0.89%	0.056%
39	15.275	2082	2089	2095	rVV	3959927	6047911	86.43%	5.408%
40	15.334	2095	2099	2106	rVB2	88169	140976	2.01%	0.126%
41	15.410	2106	2112	2119	rBV	1005163	1432445	20.47%	1.281%
42	15.516	2123	2130	2131	rVV6	24417	55874	0.80%	0.050%
43	15.628	2144	2149	2152	rBV2	44980	75590	1.08%	0.068%
44	15.775	2170	2174	2181	rVB2	30125	62807	0.90%	0.056%
45	15.845	2181	2186	2192	rBV8	19406	42655	0.61%	0.038%
46	16.010	2208	2214	2220	rVB9	39910	88654	1.27%	0.079%
47	16.181	2238	2243	2253	rVB2	108391	168339	2.41%	0.151%
48	16.310	2259	2265	2267	rBV2	24511	46213	0.66%	0.041%
49	16.434	2282	2286	2292	rBV3	77326	126887	1.81%	0.113%
50	16.563	2301	2308	2312	rBV6	31413	73965	1.06%	0.066%
51	16.687	2324	2329	2335	rVB	66410	107773	1.54%	0.096%
52	16.757	2337	2341	2343	rBV4	31317	46016	0.66%	0.041%
53	16.792	2343	2347	2352	rVB5	44485	71016	1.01%	0.064%
54	16.845	2352	2356	2362	rVB	74061	112093	1.60%	0.100%
55	16.916	2362	2368	2371	rBV2	59024	120183	1.72%	0.107%
56	17.028	2380	2387	2392	rBV	4160430	6509614	93.03%	5.821%
57	17.075	2392	2395	2399	rVV	1355342	1784945	25.51%	1.596%
58	17.128	2399	2404	2415	rVV2	4147573	6304487	90.10%	5.638%
59	17.210	2415	2418	2425	rVV2	44334	89412	1.28%	0.080%
60	17.322	2434	2437	2445	rVB2	32347	61843	0.88%	0.055%
61	17.439	2451	2457	2468	rVB	136977	250584	3.58%	0.224%
62	17.551	2469	2476	2482	rVB3	39738	83348	1.19%	0.075%
63	17.610	2482	2486	2494	rBV5	42628	93742	1.34%	0.084%
64	17.704	2498	2502	2506	rBV	94486	119061	1.70%	0.106%
65	17.745	2506	2509	2514	rVB3	32133	49661	0.71%	0.044%
66	17.810	2515	2520	2526	rBV7	56084	121296	1.73%	0.108%
67	17.904	2527	2536	2537	rBV	188173	297758	4.26%	0.266%
68	17.934	2537	2541	2550	rVB2	1594254	2454374	35.08%	2.195%
69	18.098	2563	2569	2573	rBV2	302056	510221	7.29%	0.456%
70	18.134	2573	2575	2579	rVB	102374	116148	1.66%	0.104%
71	18.263	2594	2597	2600	rVB2	49742	62245	0.89%	0.056%
72	18.410	2618	2622	2626	rBV	166674	223361	3.19%	0.200%
73	18.457	2626	2630	2635	rVB	232510	307815	4.40%	0.275%
74	18.651	2661	2663	2668	rVB3	60783	81509	1.16%	0.073%
75	18.828	2690	2693	2696	rBV2	73126	102182	1.46%	0.091%
76	18.922	2707	2709	2713	rVB2	63115	68843	0.98%	0.062%

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Stop Thrs : 0

Peak Location: TOP

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77	18.975	2713	2718	2725	rBV3	137901	323656	4.63%	0.289%
78	19.092	2730	2738	2746	rBV	3183689	4669949	66.74%	4.176%
79	19.222	2756	2760	2768	rVB2	547735	834384	11.92%	0.746%
80	19.428	2790	2795	2798	rBV	4441105	6611083	94.48%	5.912%
81	19.457	2798	2800	2808	rVV	2125580	2647000	37.83%	2.367%
82	19.533	2809	2813	2818	rVB2	113331	164666	2.35%	0.147%
83	19.781	2851	2855	2862	rBV2	143620	331928	4.74%	0.297%
84	19.951	2881	2884	2887	rBV	272609	365611	5.22%	0.327%
85	20.051	2898	2901	2905	rBV2	232927	306141	4.38%	0.274%
86	20.363	2952	2954	2958	rVB	187247	187613	2.68%	0.168%
87	20.469	2969	2972	2976	rVB	326323	398508	5.70%	0.356%
88	20.704	3009	3012	3018	rVB	216232	286491	4.09%	0.256%
89	20.916	3045	3048	3050	rBV	163596	209548	2.99%	0.187%
90	20.945	3050	3053	3060	rVB3	1486722	2163427	30.92%	1.935%
91	21.169	3087	3091	3095	rVB	558982	674996	9.65%	0.604%
92	21.263	3099	3107	3111	rVV3	3646309	6997345	100.00%	6.257%
93	21.304	3111	3114	3125	rVB	1091216	1675449	23.94%	1.498%
94	21.445	3135	3138	3145	rVB2	406695	578517	8.27%	0.517%
95	21.998	3227	3232	3234	rBV2	642503	1012916	14.48%	0.906%
96	23.257	3439	3446	3452	rBV	705618	2003320	28.63%	1.791%
97	23.333	3454	3459	3466	rVB2	1048032	2207617	31.55%	1.974%
98	23.974	3561	3568	3574	rBV2	1108688	2598065	37.13%	2.323%
99	24.174	3594	3602	3612	rVB	2036221	4978594	71.15%	4.452%
100	25.168	3766	3771	3783	rVB2	462426	1285554	18.37%	1.150%

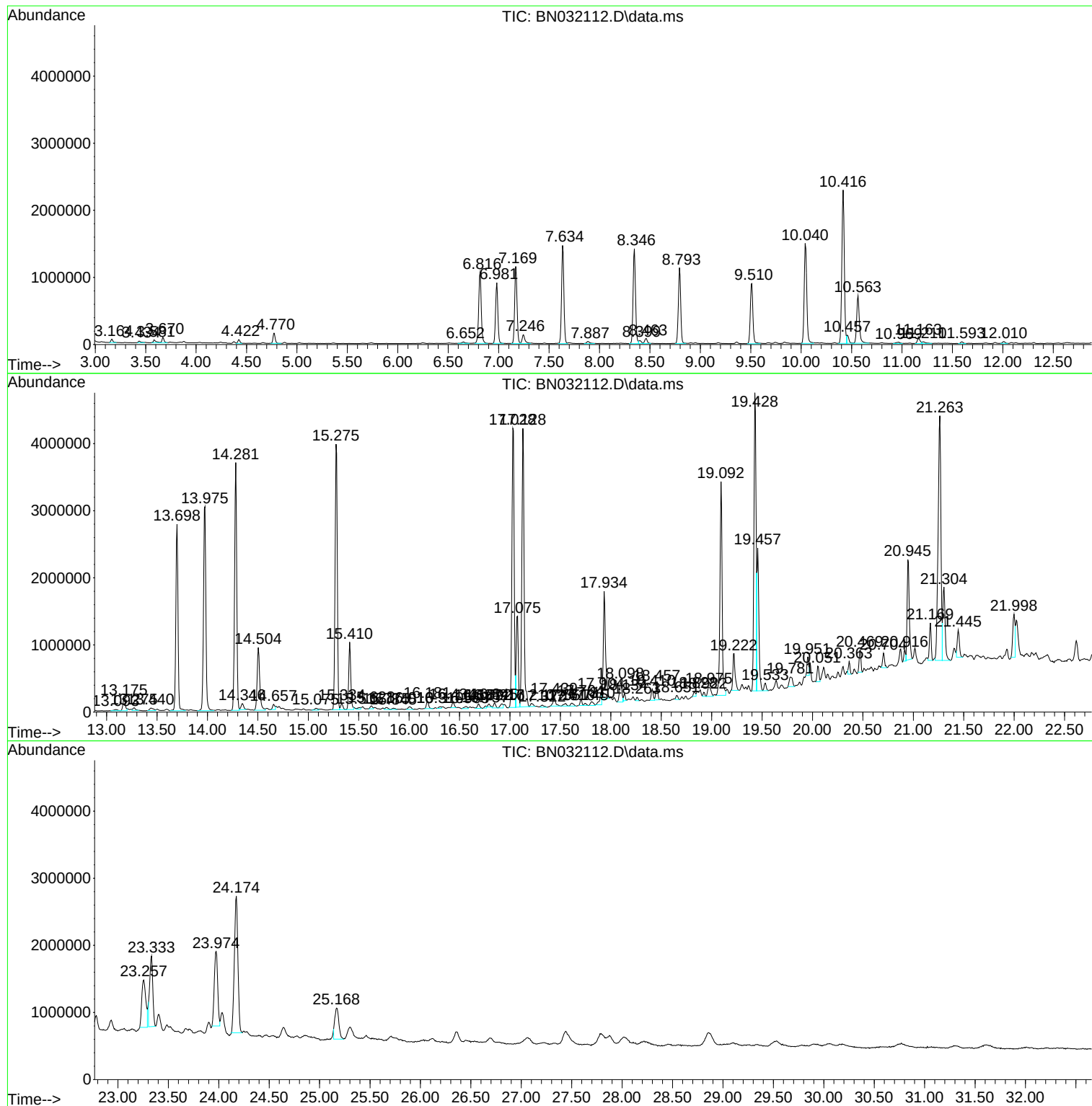
Sum of corrected areas: 111827837

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

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



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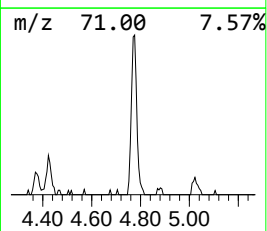
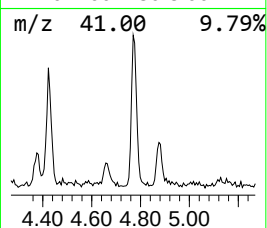
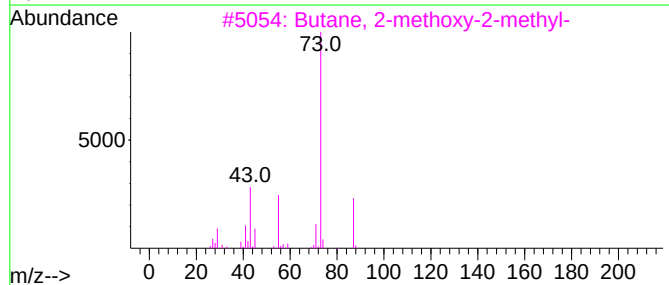
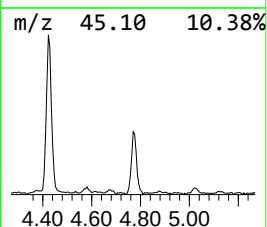
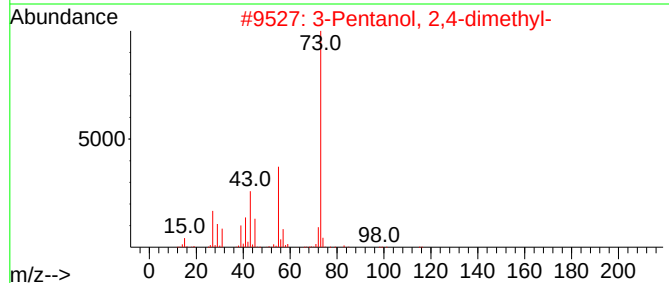
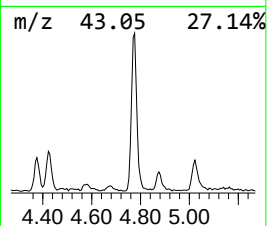
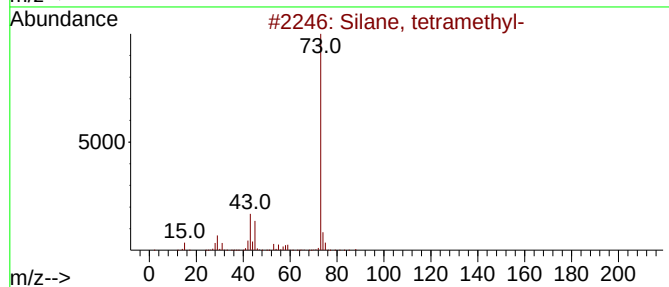
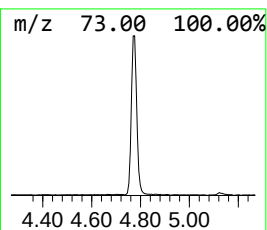
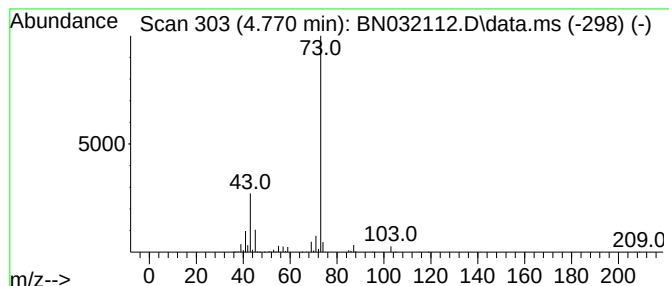
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN052924.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 unknown-01 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.770	2.02 ng/ul	237432	1,4-Dichlorobenzene-d4	7.634

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Silane, tetramethyl-	88	C4H12Si	000075-76-3	38
2			3-Pentanol, 2,4-dimethyl-	116	C7H16O	000600-36-2	36
3			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	36
4			Octanal, 7-methoxy-3,7-dimethyl-	186	C11H22O2	003613-30-7	33
5			Acetic acid, 3-[1,3]dioxolan-2-y...	174	C8H14O4	1000186-02-3	33



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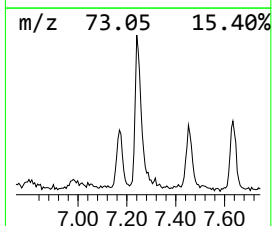
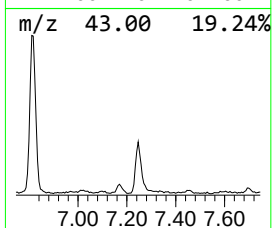
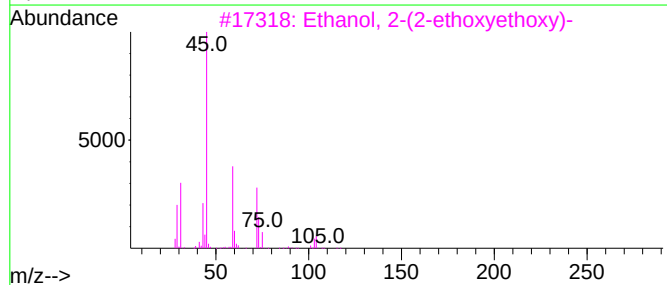
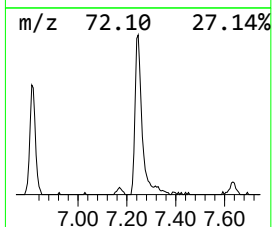
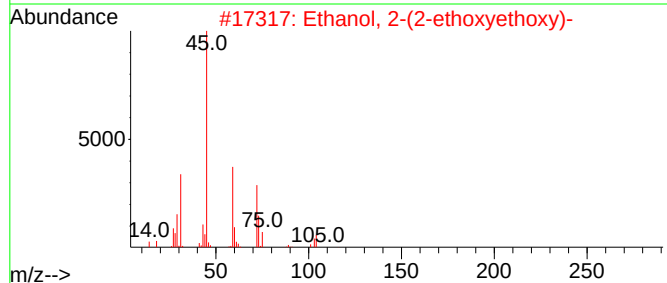
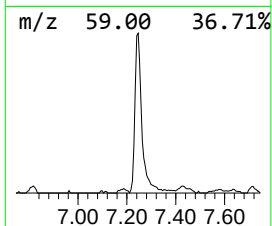
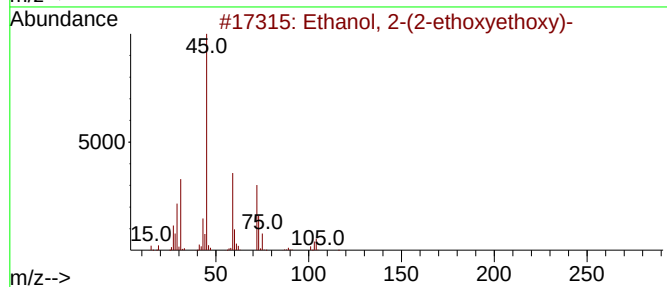
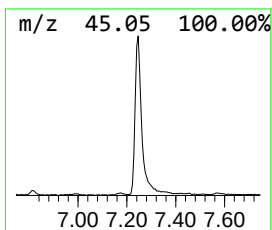
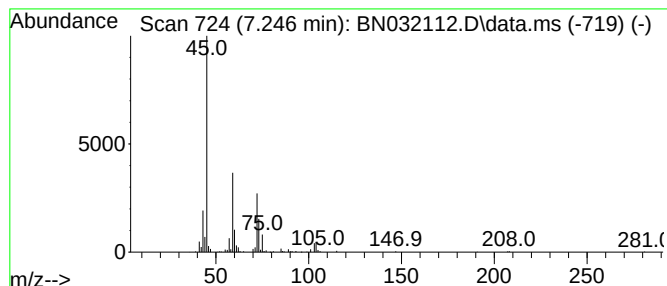
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN052924.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Ethanol, 2-(2-ethoxyethoxy)- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.246	2.16 ng/ul	254267	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	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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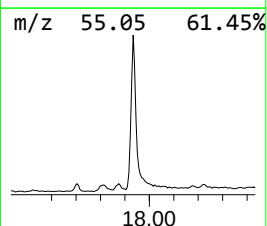
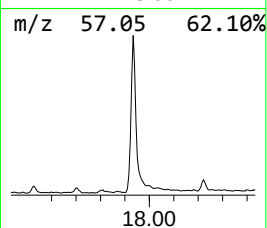
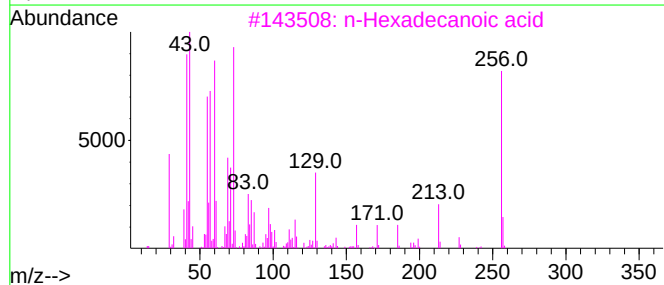
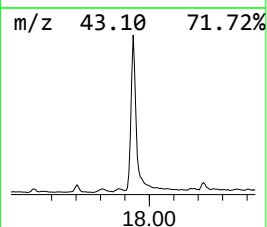
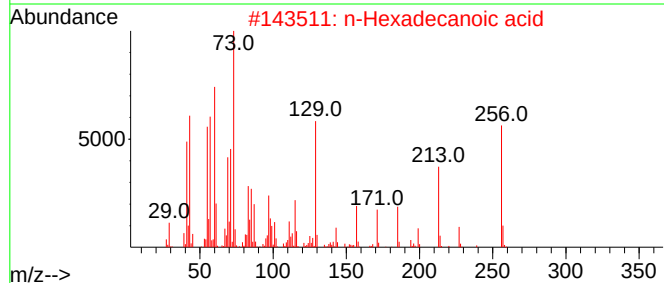
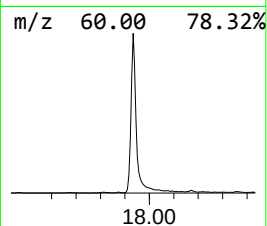
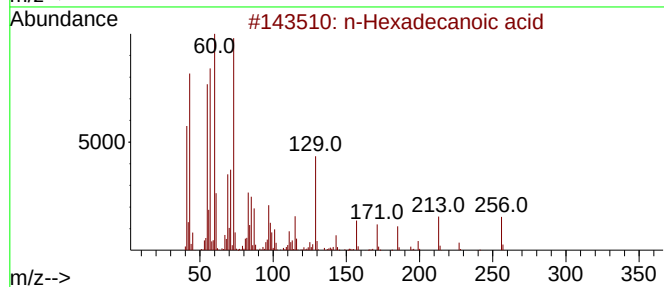
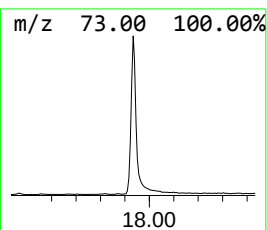
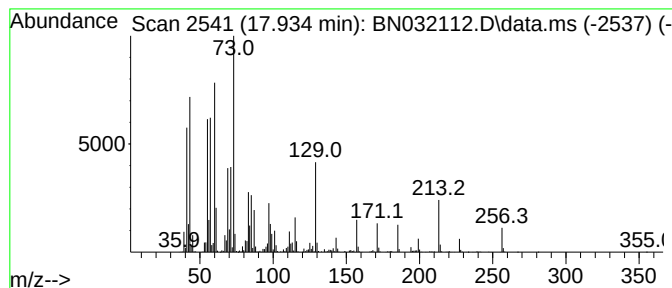
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN052924.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 n-Hexadecanoic acid Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.934	7.54 ng/ul	2454370	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	98
5			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062024\
Data File : BN032112.D
Acq On : 21 Jun 2024 01:14
Operator : MA/JU
Sample : P2710-07
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
A4BL4

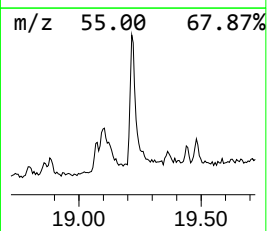
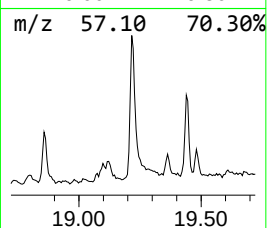
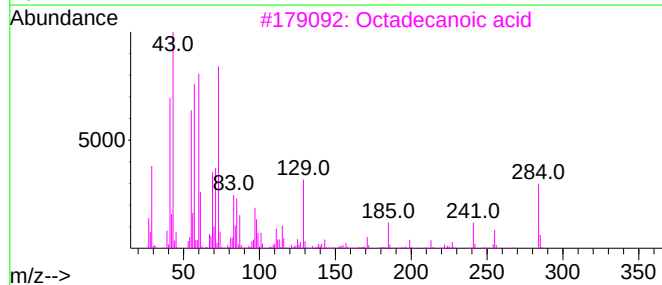
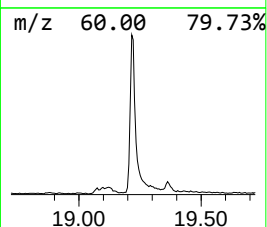
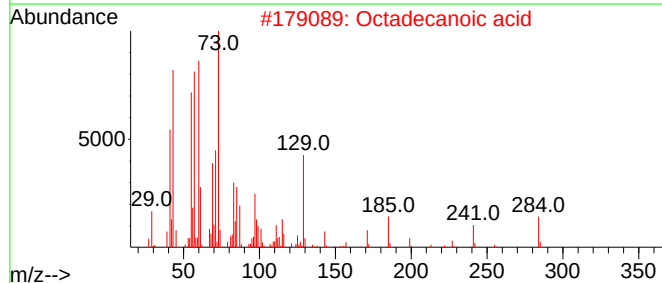
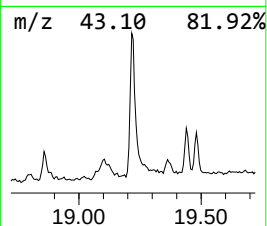
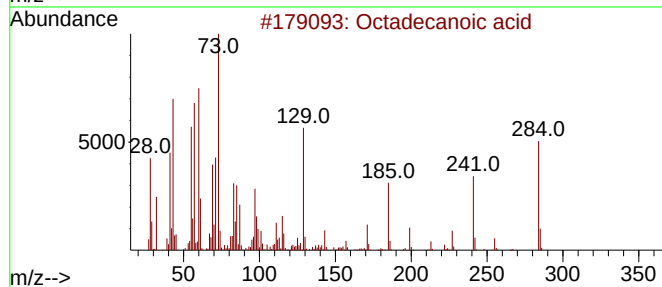
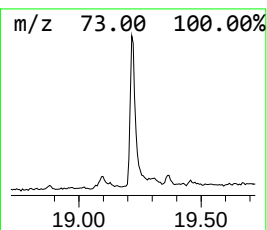
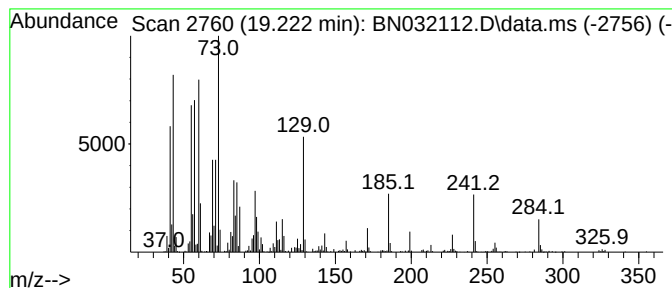
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Octadecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.222	2.38 ng/ul	834384	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	97
3			Octadecanoic acid	284	C18H36O2	000057-11-4	95
4			Octadecanoic acid	284	C18H36O2	000057-11-4	95
5			Octadecanoic acid, 2-(2-hydroxye...	372	C22H44O4	000106-11-6	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062024\
Data File : BN032112.D
Acq On : 21 Jun 2024 01:14
Operator : MA/JU
Sample : P2710-07
Misc :
ALS Vial : 20 Sample Multiplier: 1

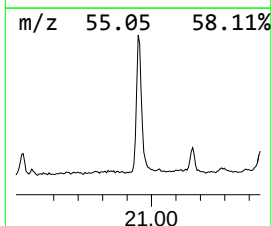
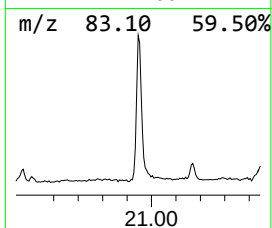
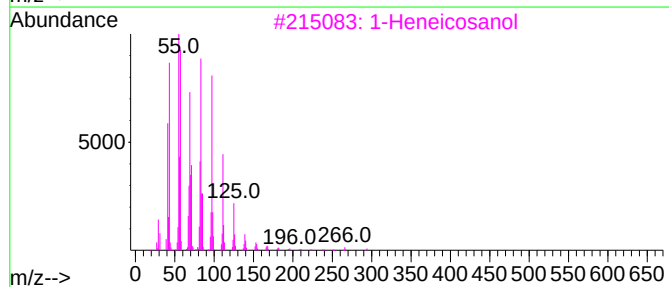
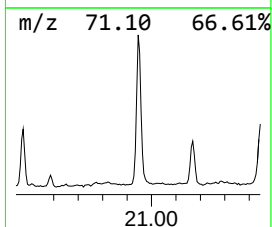
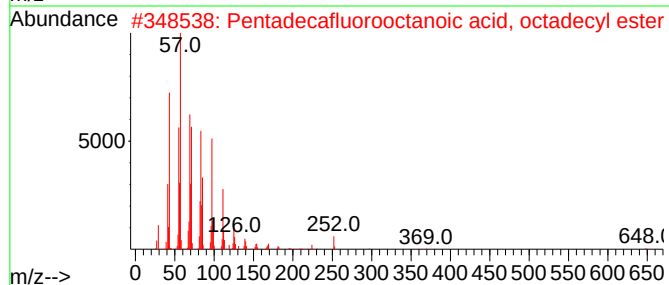
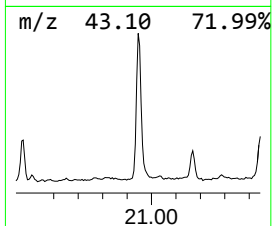
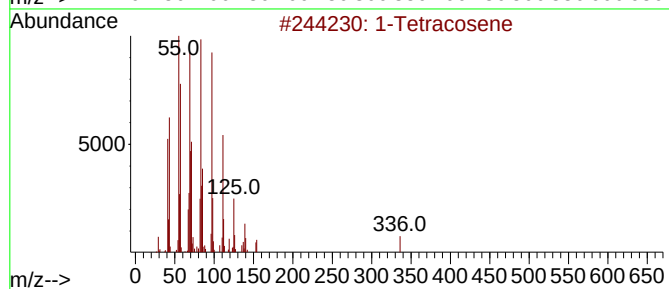
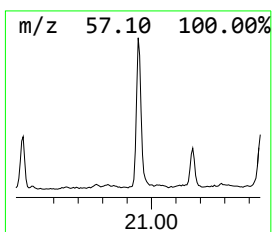
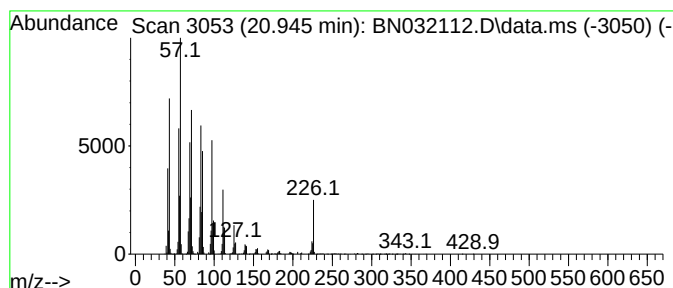
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN052924.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
20.945	6.18 ng/ul	2163430	Chrysene-d12	21.263	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Tetracosene	336	C24H48	010192-32-2	93
2	Pentadecafluorooctanoic acid, oc...	666	C26H37F15O2	1000406-04-8	93
3	1-Heneicosanol	312	C21H44O	015594-90-8	91
4	Cyclohexadecane	224	C16H32	000295-65-8	91
5	Ethanol, 2-(tetradecyloxy)-	258	C16H34O2	002136-70-1	76



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062024\
Data File : BN032112.D
Acq On : 21 Jun 2024 01:14
Operator : MA/JU
Sample : P2710-07
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
A4BL4

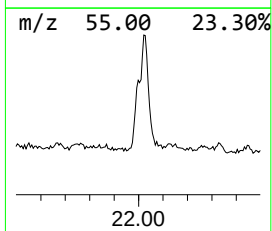
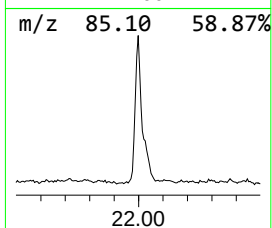
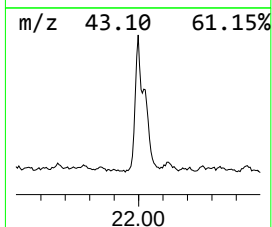
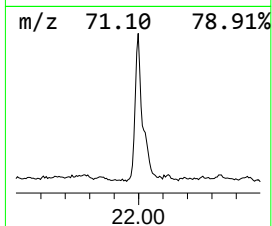
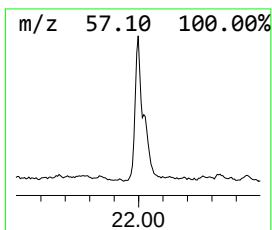
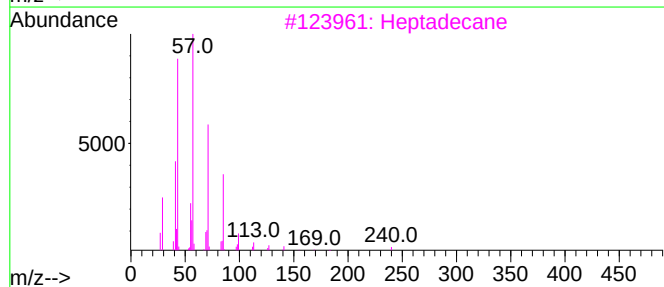
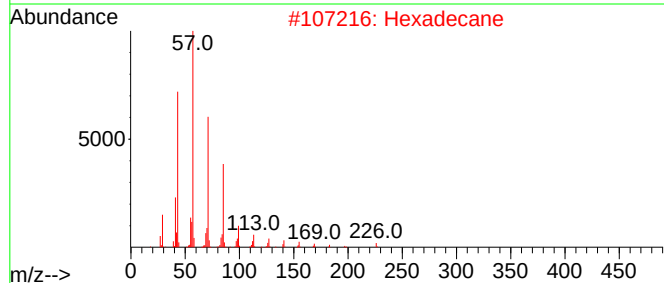
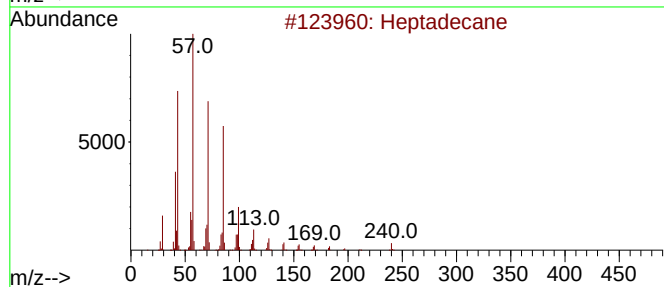
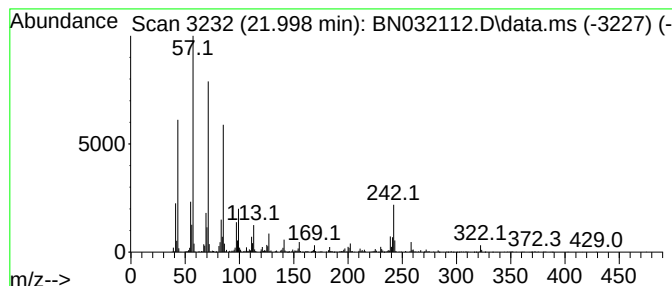
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.998	2.90 ng/ul	1012920	Chrysene-d12	21.263

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heptadecane	240	C17H36	000629-78-7	96
2		Hexadecane	226	C16H34	000544-76-3	95
3		Heptadecane	240	C17H36	000629-78-7	95
4		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	94
5		Heptacosane	380	C27H56	000593-49-7	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062024\
Data File : BN032112.D
Acq On : 21 Jun 2024 01:14
Operator : MA/JU
Sample : P2710-07
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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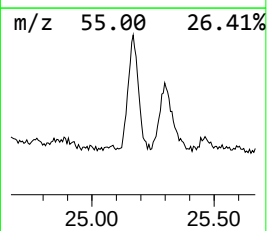
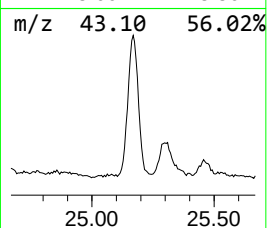
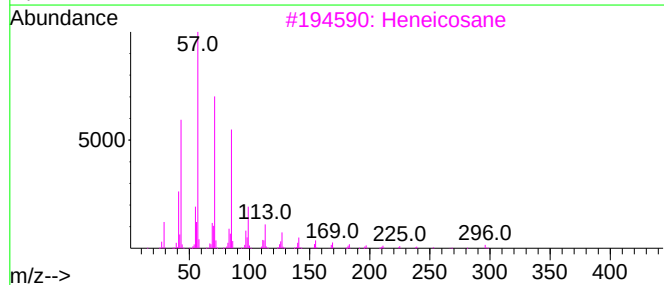
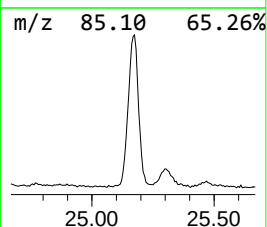
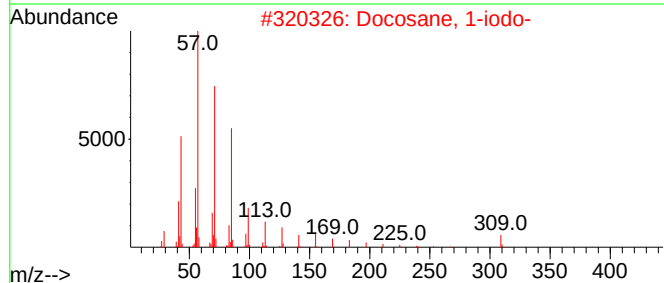
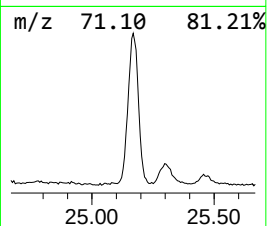
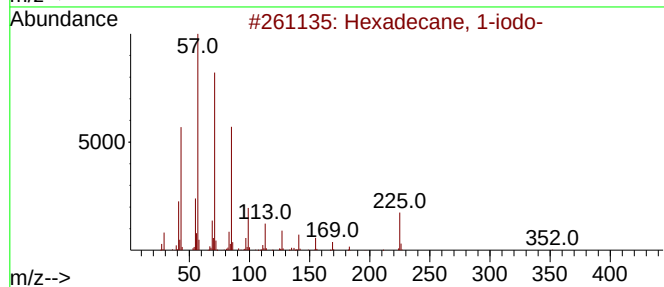
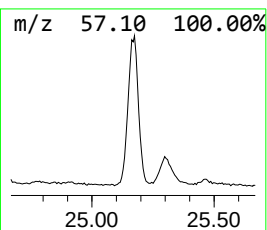
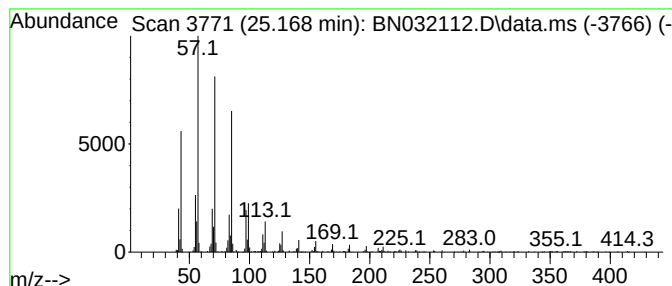
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Hexadecane, 1-iodo- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.169	5.16 ng/ul	1285550	Perylene-d12	24.174

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexadecane, 1-iodo-	352	C16H33I	000544-77-4	93
2			Docosane, 1-iodo-	436	C22H45I	1000406-31-9	90
3			Heneicosane	296	C21H44	000629-94-7	90
4			Octacosane, 2-methyl-	408	C29H60	001560-98-1	90
5			Octadecane, 1-iodo-	380	C18H37I	000629-93-6	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062024\
Data File : BN032112.D
Acq On : 21 Jun 2024 01:14
Operator : MA/JU
Sample : P2710-07
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
A4BL4

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
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Ethanol, 2-(2-e...	7.246	2.2	ng/u1	254267	1	7.634	2352180	20.0
n-Hexadecanoic ...	17.934	7.5	ng/u1	2454370	4	17.028	6509610	20.0
Octadecanoic acid	19.222	2.4	ng/u1	834384	5	21.263	6997350	20.0
(DEL) Alkane: S...	20.945	6.2	ng/u1	2163430	5	21.263	6997350	20.0
(DEL) Alkane: S...	21.998	2.9	ng/u1	1012920	5	21.263	6997350	20.0
Hexadecane, 1-i...	25.169	5.2	ng/u1	1285550	6	24.174	4978590	20.0