

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN122022\
 Data File : BN023296.D
 Acq On : 20 Dec 2022 13:38
 Operator : CG/JU
 Sample : N6003-18
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 DBXT8

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-BN121522.M
 Title : SVOA CALIBRATION

Signal : TIC: BN023296.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.340	39	43	51	rVB	76699	113072	1.13%	0.088%
2	3.770	114	116	132	rBV	572603	968012	9.66%	0.752%
3	3.887	133	136	142	rVB	27395	37534	0.37%	0.029%
4	4.017	153	158	164	rVB2	25863	42682	0.43%	0.033%
5	4.111	171	174	185	rVB	20662	35752	0.36%	0.028%
6	5.064	329	336	345	rVB	263096	388654	3.88%	0.302%
7	7.152	683	691	702	rBV	1542612	2369255	23.64%	1.840%
8	7.322	713	720	730	rVB	1286034	1936202	19.32%	1.504%
9	7.522	744	754	765	rBV	1616614	2485551	24.80%	1.931%
10	7.611	765	769	777	rVB4	10369	18751	0.19%	0.015%
11	7.993	825	834	842	rBV	1043457	1669282	16.66%	1.297%
12	8.710	945	956	967	rBV	1720651	2737137	27.31%	2.126%
13	8.828	971	976	981	rVB	33153	52158	0.52%	0.041%
14	9.169	1025	1034	1044	rBV	1543987	2530209	25.25%	1.965%
15	9.581	1095	1104	1114	rVB2	25510	48793	0.49%	0.038%
16	9.763	1131	1135	1140	rVB	16556	25463	0.25%	0.020%
17	9.893	1149	1157	1170	rBV	1354467	2189573	21.85%	1.701%
18	10.116	1191	1195	1201	rVB5	12074	19799	0.20%	0.015%
19	10.434	1240	1249	1259	rBV	2048006	3372985	33.65%	2.620%
20	10.593	1270	1276	1284	rVB2	27106	47065	0.47%	0.037%
21	10.816	1306	1314	1320	rBV	1609978	2631260	26.25%	2.044%
22	10.863	1320	1322	1326	rVB	14553	19434	0.19%	0.015%
23	10.952	1328	1337	1356	rBV	1616921	2793709	27.87%	2.170%
24	11.599	1442	1447	1451	rBV6	10403	18409	0.18%	0.014%
25	12.234	1548	1555	1560	rBV	21906	33861	0.34%	0.026%
26	12.398	1578	1583	1590	rVB	38275	59691	0.60%	0.046%
27	13.363	1738	1747	1749	rBV7	9723	17315	0.17%	0.013%
28	13.469	1761	1765	1771	rVB	43204	58157	0.58%	0.045%
29	13.675	1797	1800	1809	rVB4	12429	21617	0.22%	0.017%
30	13.969	1846	1850	1854	rBV6	14782	22098	0.22%	0.017%
31	14.057	1858	1865	1874	rVV	3762825	5005713	49.95%	3.888%
32	14.169	1880	1884	1889	rVB6	13516	21932	0.22%	0.017%
33	14.240	1892	1896	1900	rVB	15417	19683	0.20%	0.015%
34	14.340	1906	1913	1925	rVB	4350311	6314907	63.01%	4.905%
35	14.540	1938	1947	1952	rVB	26096	45600	0.45%	0.035%
36	14.645	1956	1965	1973	rBV	2604297	3709520	37.01%	2.881%

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Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN121522.M
 Title : SVOA CALIBRATION

37	14.716	1973	1977	1983	rVB8	13793	24820	0.25%	0.019%
38	14.840	1987	1998	2016	rBV	1773180	2573114	25.67%	1.999%
39	14.992	2020	2024	2029	rVV2	24947	41149	0.41%	0.032%
40	15.057	2030	2035	2041	rVB4	53118	87443	0.87%	0.068%
41	15.204	2056	2060	2068	rBV3	11811	23891	0.24%	0.019%
42	15.639	2126	2134	2140	rBV	5487552	7728263	77.11%	6.003%
43	15.692	2140	2143	2147	rVV5	15538	25798	0.26%	0.020%
44	15.751	2147	2153	2166	rVV	1780818	2407644	24.02%	1.870%
45	15.875	2169	2174	2183	rVB	31178	51036	0.51%	0.040%
46	16.004	2190	2196	2203	rVB	1806603	2361121	23.56%	1.834%
47	16.204	2224	2230	2236	rVB10	8917	20217	0.20%	0.016%
48	16.363	2251	2257	2261	rVV8	13528	31785	0.32%	0.025%
49	16.563	2287	2291	2297	rVB	44521	63388	0.63%	0.049%
50	16.786	2325	2329	2336	rVB8	13926	23272	0.23%	0.018%
51	17.092	2375	2381	2385	rVB3	23510	41048	0.41%	0.032%
52	17.145	2387	2390	2393	rBV5	14317	16838	0.17%	0.013%
53	17.286	2410	2414	2415	rBV4	14419	17833	0.18%	0.014%
54	17.392	2426	2432	2437	rBV	3287387	4548001	45.38%	3.533%
55	17.434	2437	2439	2443	rVV	88686	99633	0.99%	0.077%
56	17.492	2443	2449	2458	rVV	5733026	8100440	80.82%	6.292%
57	17.581	2460	2464	2473	rVB5	56840	96060	0.96%	0.075%
58	17.675	2475	2480	2482	rVV4	14252	25421	0.25%	0.020%
59	17.704	2482	2485	2489	rVV3	27150	38355	0.38%	0.030%
60	17.769	2493	2496	2504	rVB5	32390	58520	0.58%	0.045%
61	18.016	2534	2538	2541	rBV3	12987	17062	0.17%	0.013%
62	18.057	2542	2545	2550	rVB3	31606	40259	0.40%	0.031%
63	18.110	2550	2554	2557	rBV5	12738	18328	0.18%	0.014%
64	18.169	2558	2564	2569	rBV3	36660	68225	0.68%	0.053%
65	18.228	2569	2574	2579	rBV	117825	187685	1.87%	0.146%
66	18.286	2579	2584	2594	rBV	1152044	1699808	16.96%	1.320%
67	18.481	2613	2617	2622	rVB2	68881	103603	1.03%	0.080%
68	18.539	2622	2627	2631	rVB	83370	107730	1.07%	0.084%
69	18.616	2637	2640	2645	rVB3	27005	37213	0.37%	0.029%
70	18.710	2651	2656	2663	rBV2	323772	396300	3.95%	0.308%
71	18.775	2663	2667	2668	rBV3	29962	37985	0.38%	0.030%
72	18.804	2668	2672	2677	rVB5	57075	80173	0.80%	0.062%
73	18.869	2678	2683	2688	rBV	371021	486895	4.86%	0.378%
74	19.139	2726	2729	2733	rBV3	38207	59920	0.60%	0.047%
75	19.216	2734	2742	2746	rVV	235098	337493	3.37%	0.262%
76	19.251	2746	2748	2753	rVB3	42621	51298	0.51%	0.040%

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 Title : SVOA CALIBRATION

77	19.304	2754	2757	2761	rBV2	32422	63145	0.63%	0.049%
78	19.351	2761	2765	2769	rVB	105454	132595	1.32%	0.103%
79	19.416	2772	2776	2778	rBV	171518	213980	2.14%	0.166%
80	19.445	2778	2781	2795	rVB	345539	617926	6.17%	0.480%
81	19.563	2796	2801	2809	rBV	435074	687426	6.86%	0.534%
82	19.786	2833	2839	2849	rBV2	7219008	10022392	100.00%	7.785%
83	19.857	2849	2851	2859	rVB7	39993	60858	0.61%	0.047%
84	19.945	2861	2866	2871	rBV	2936161	3435394	34.28%	2.668%
85	20.069	2884	2887	2893	rVV2	69592	103317	1.03%	0.080%
86	20.127	2894	2897	2905	rVB5	87965	155634	1.55%	0.121%
87	20.298	2922	2926	2933	rVB3	324200	475484	4.74%	0.369%
88	20.539	2964	2967	2975	rVB3	113907	155717	1.55%	0.121%
89	21.274	3077	3092	3102	rBV3	1862871	4586346	45.76%	3.562%
90	21.522	3131	3134	3137	rBV3	176887	268352	2.68%	0.208%
91	21.580	3138	3144	3148	rBV	3847217	5717855	57.05%	4.441%
92	21.727	3167	3169	3178	rVB4	127700	153636	1.53%	0.119%
93	22.198	3245	3249	3251	rBV2	178461	267562	2.67%	0.208%
94	23.269	3426	3431	3436	rBV3	561209	994815	9.93%	0.773%
95	23.868	3526	3533	3548	rVB2	4141827	8900308	88.80%	6.913%
96	24.027	3553	3560	3575	rVB2	2149183	4495084	44.85%	3.491%
97	24.651	3660	3666	3675	rVB	716282	1494160	14.91%	1.161%
98	24.751	3678	3683	3696	rVB2	347567	854579	8.53%	0.664%
99	26.545	3980	3988	4003	rVB3	1005390	3259089	32.52%	2.531%
100	29.092	4409	4421	4433	rBV5	1795902	7514078	74.97%	5.836%

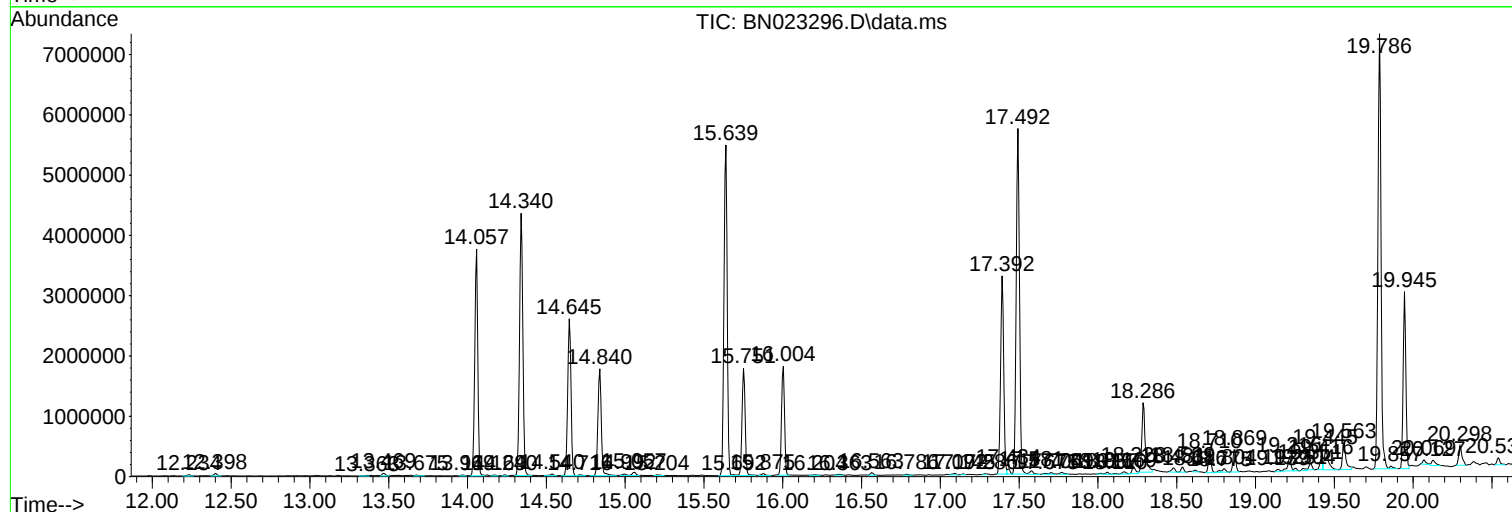
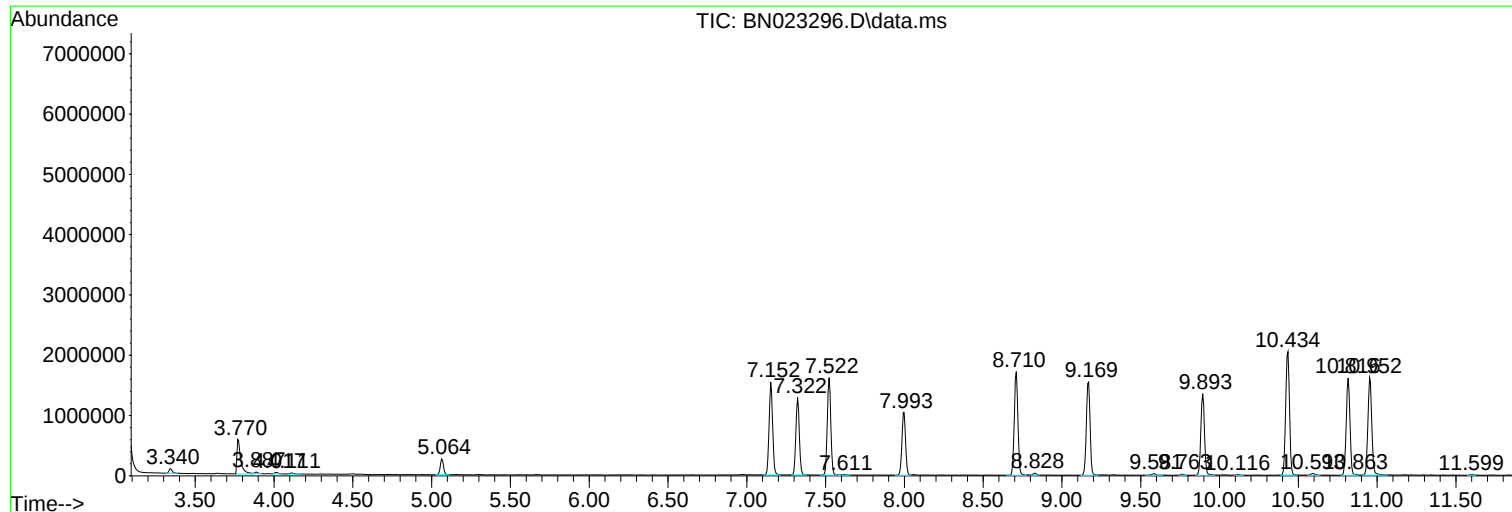
Sum of corrected areas: 128745634

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Instrument :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN121522.M
Quant Title : SVOA CALIBRATION

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



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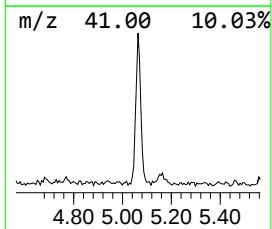
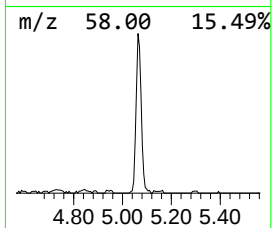
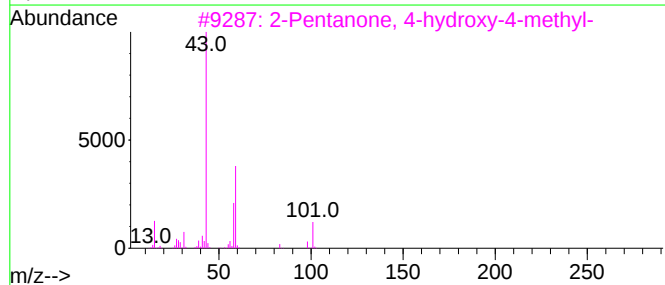
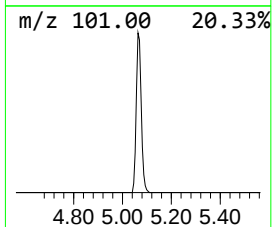
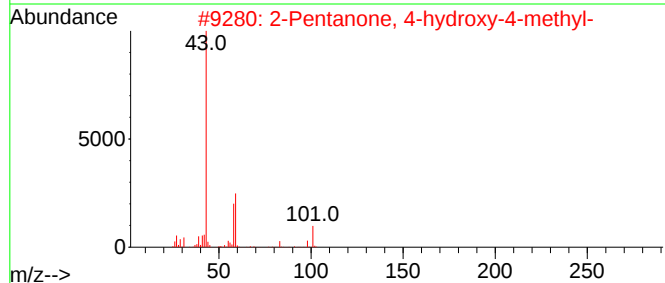
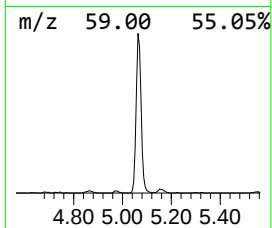
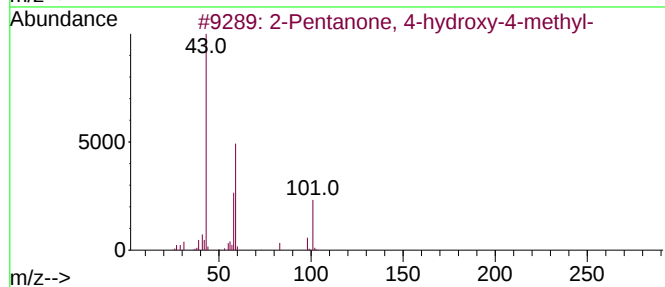
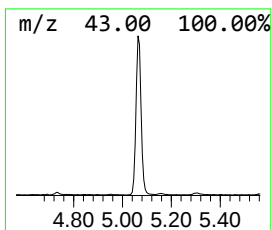
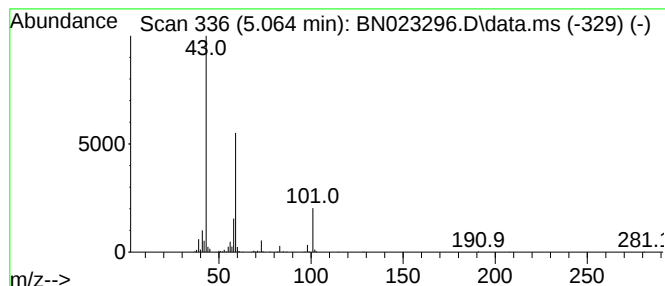
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN121522.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.064	4.66 ng/ul	388654	1,4-Dichlorobenzene-d4	7.993

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56		
2	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50		
3	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	45		
4	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	40		
5	Propane, 1-ethoxy-2-methyl-	102	C6H14O	000627-02-1	23		



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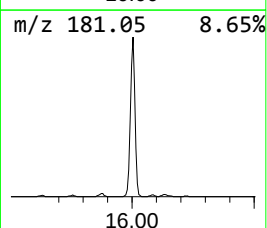
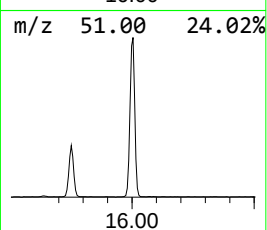
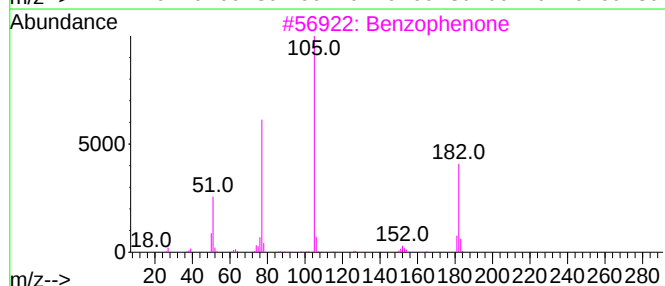
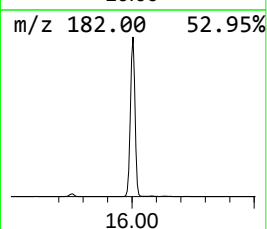
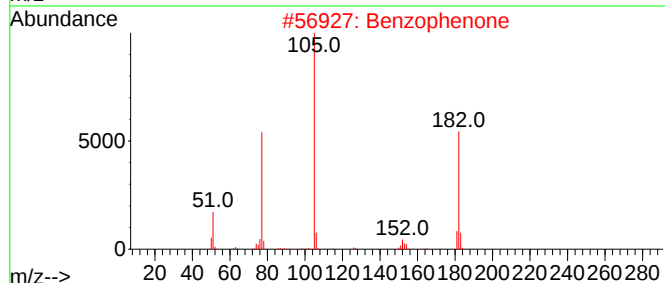
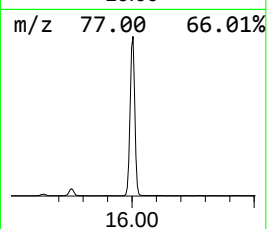
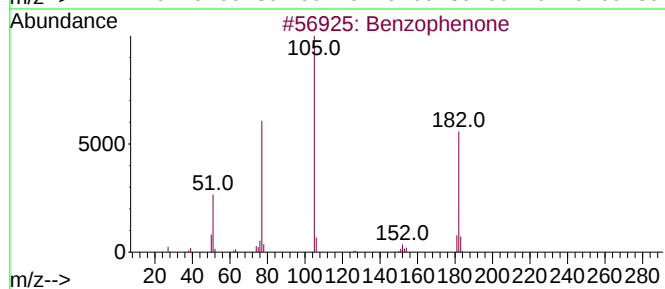
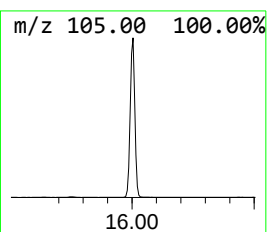
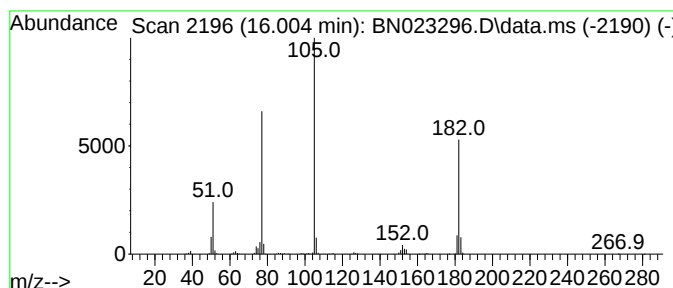
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN121522.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Benzophenone Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.004	12.73 ng/ul	2361120	Acenaphthene-d10	14.645

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	96
2			Benzophenone	182	C13H10O	000119-61-9	95
3			Benzophenone	182	C13H10O	000119-61-9	94
4			Benzophenone	182	C13H10O	000119-61-9	93
5			Benzophenone	182	C13H10O	000119-61-9	91



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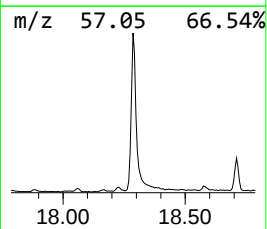
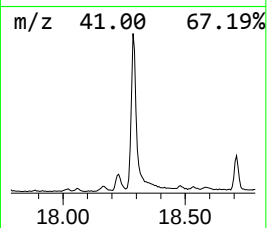
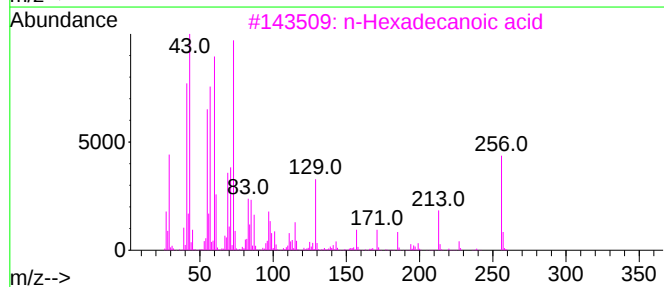
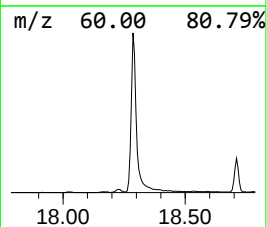
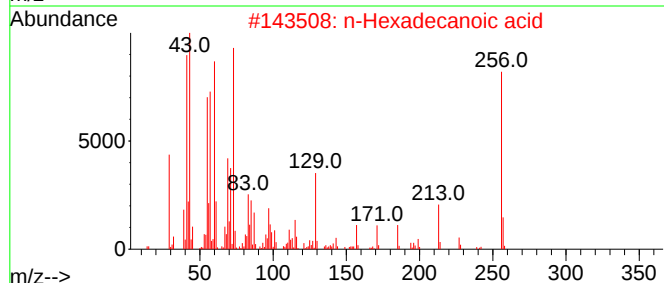
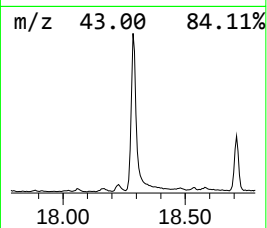
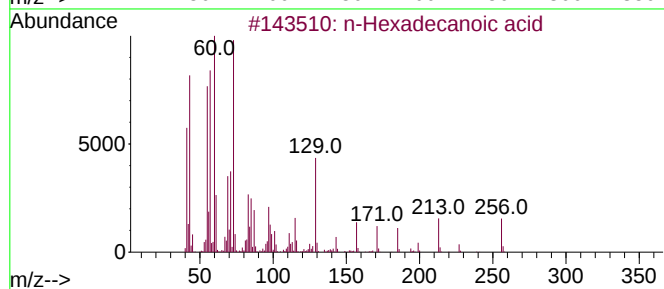
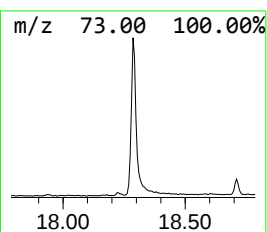
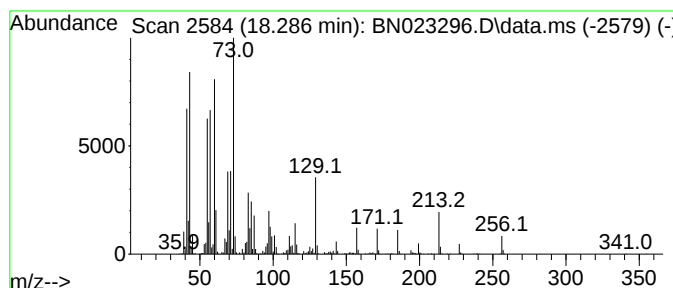
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN121522.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 n-Hexadecanoic acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.286	7.47 ng/ul	1699810	Phenanthrene-d10	17.392

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid			256	C16H32O2	000057-10-3	99
2	n-Hexadecanoic acid			256	C16H32O2	000057-10-3	99
3	n-Hexadecanoic acid			256	C16H32O2	000057-10-3	98
4	n-Hexadecanoic acid			256	C16H32O2	000057-10-3	98
5	Tetradecanoic acid			228	C14H28O2	000544-63-8	91



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Data File : BN023296.D
Acq On : 20 Dec 2022 13:38
Operator : CG/JU
Sample : N6003-18
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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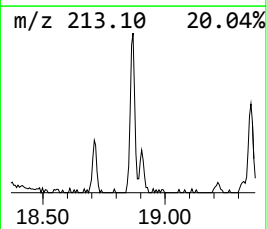
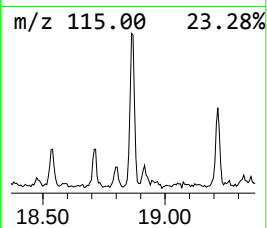
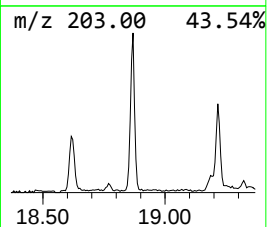
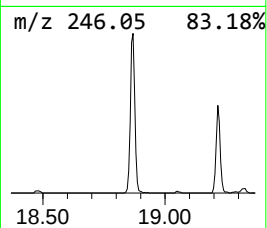
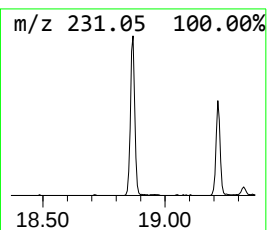
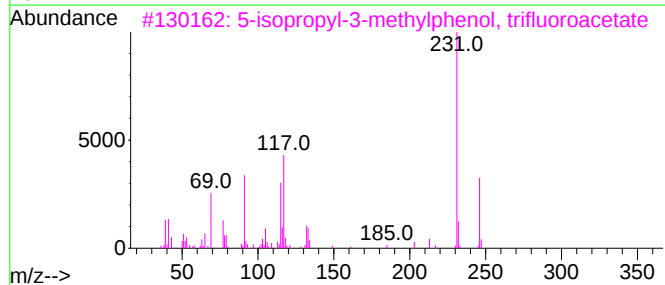
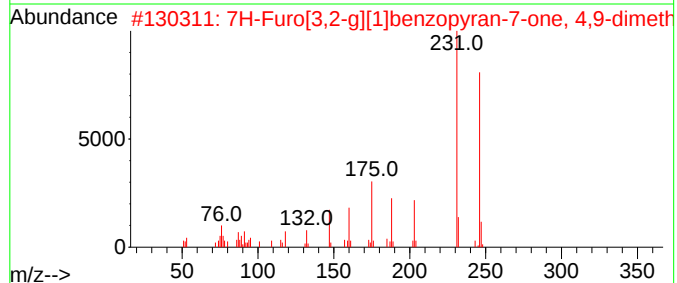
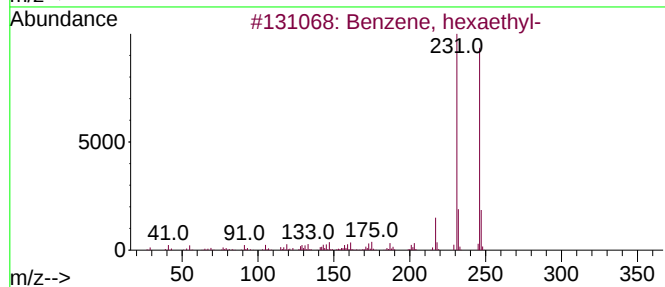
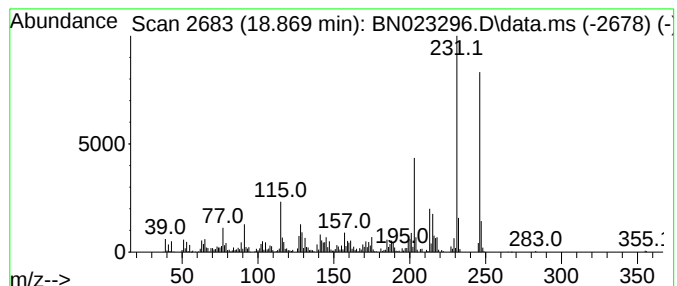
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Benzene, hexaethyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.869	2.14 ng/ul	486895	Phenanthrene-d10	17.392

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, hexaethyl-	246	C18H30	000604-88-6	86
2			7H-Furo[3,2-g][1]benzopyran-7-on...	246	C13H10O5	000482-27-9	64
3			5-isopropyl-3-methylphenol, trif...	246	C12H13F3O2	1000376-51-9	62
4			7H-Furo[3,2-g][1]benzopyran-7-on...	246	C13H10O5	000482-27-9	62
5			7H-Furo[3,2-g][1]benzopyran-7-on...	246	C13H10O5	000482-27-9	62



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN122022\
Data File : BN023296.D
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Sample : N6003-18
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Instrument :
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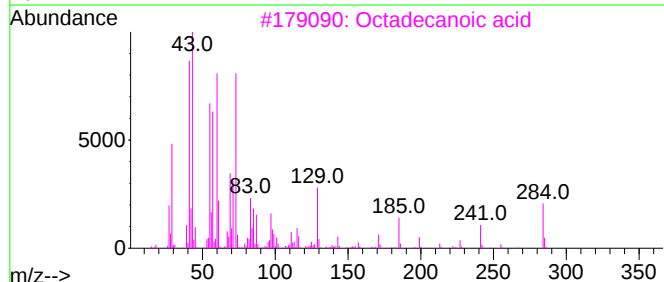
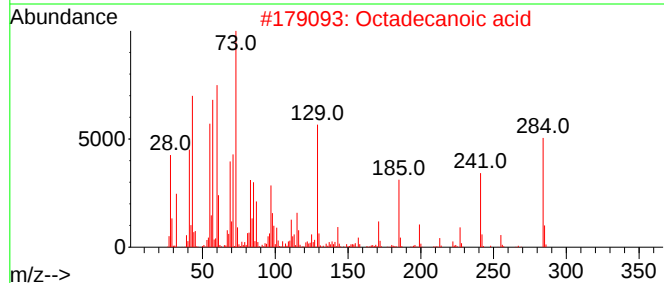
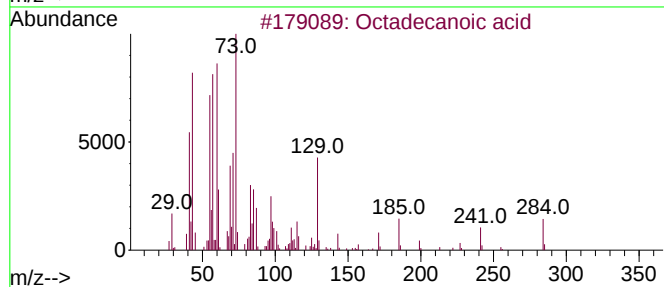
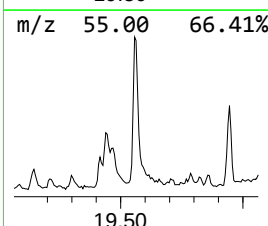
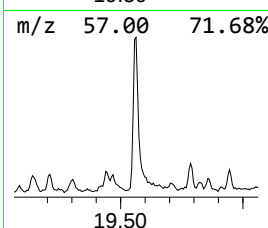
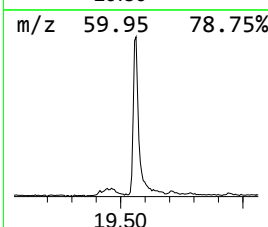
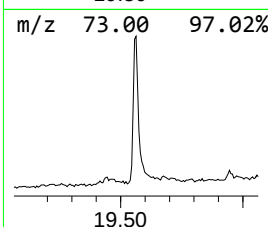
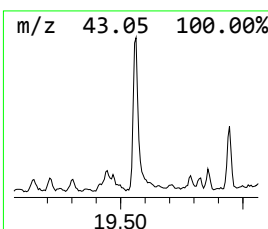
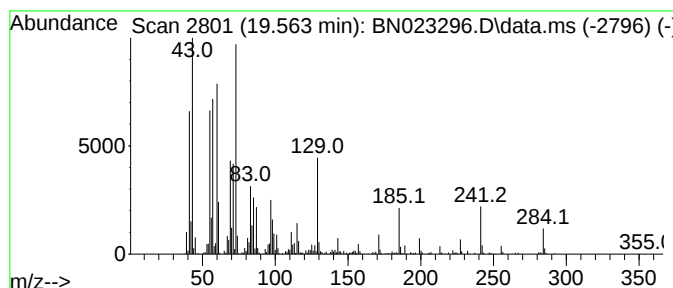
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Octadecanoic acid Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.563	2.40 ng/ul	687426	Chrysene-d12	21.580

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	98
4			Octadecanoic acid	284	C18H36O2	000057-11-4	98
5			Pentadecanoic acid	242	C15H30O2	001002-84-2	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN122022\
Data File : BN023296.D
Acq On : 20 Dec 2022 13:38
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Misc :
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Instrument :
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ClientSampleId :
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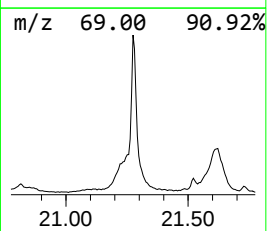
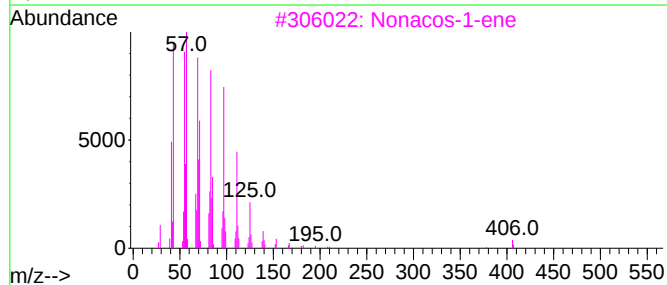
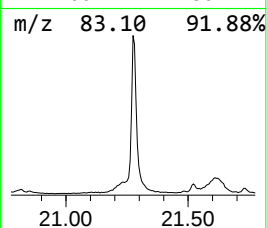
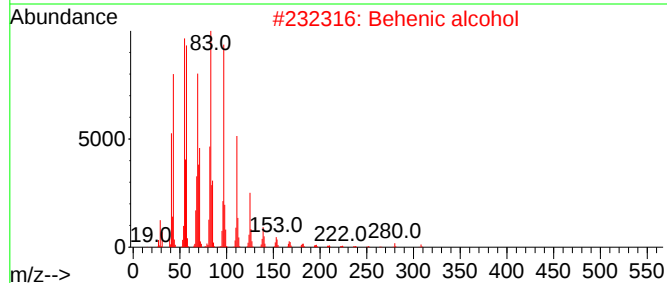
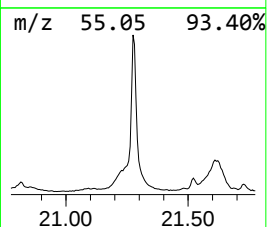
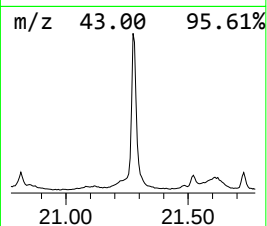
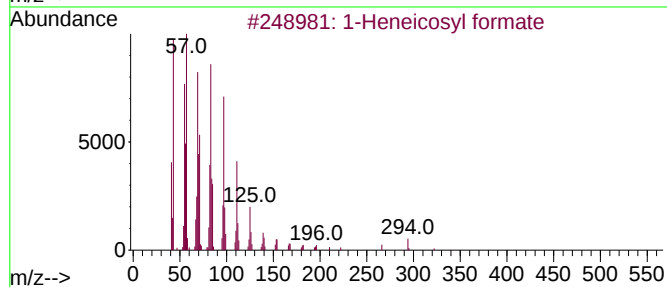
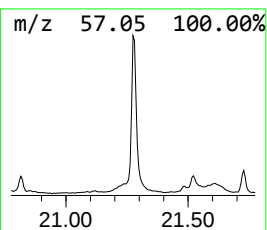
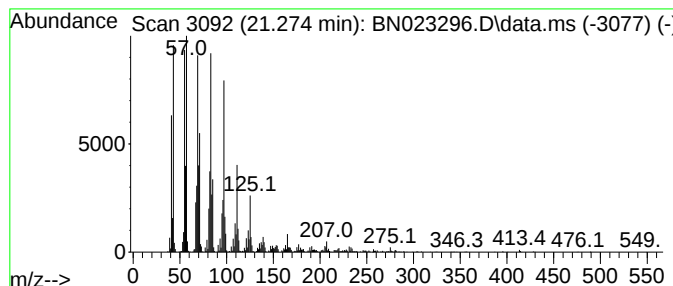
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 1-Heneicosyl formate Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.275	16.04 ng/ul	4586350	Chrysene-d12	21.580

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Heneicosyl formate	340	C22H44O2	077899-03-7	96
2			Behenic alcohol	326	C22H46O	000661-19-8	94
3			Nonacos-1-ene	406	C29H58	018835-35-3	94
4			1-Hexacosene	364	C26H52	018835-33-1	94
5			Cyclopentadecane	210	C15H30	000295-48-7	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN122022\
Data File : BN023296.D
Acq On : 20 Dec 2022 13:38
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Sample : N6003-18
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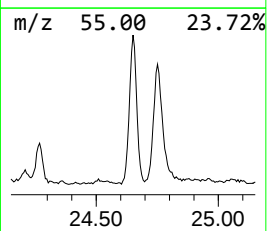
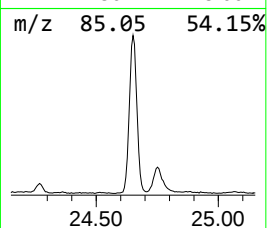
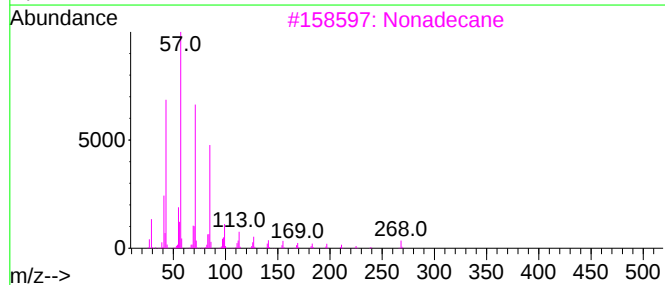
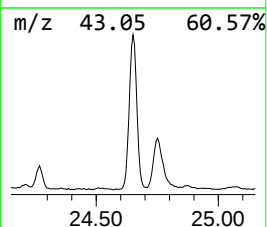
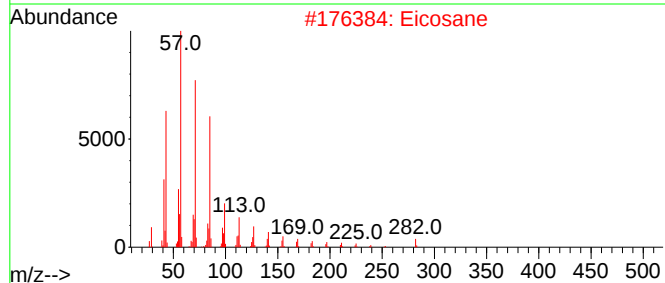
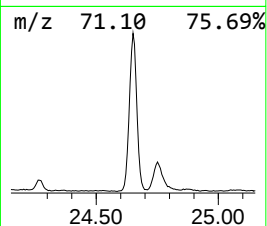
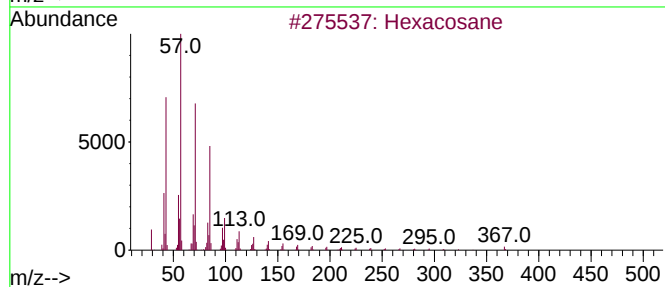
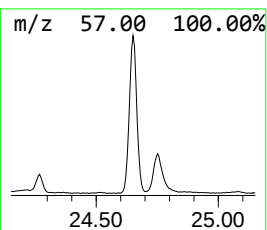
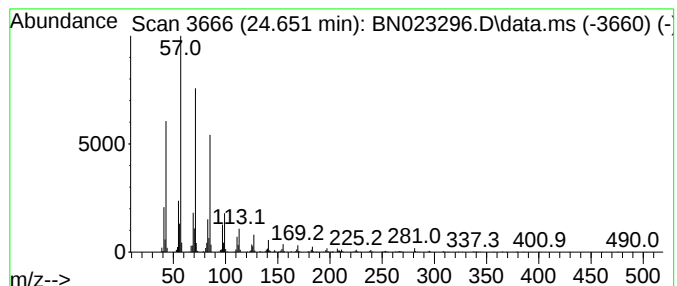
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.651	6.65 ng/ul	1494160	Perylene-d12	24.027

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Hexacosane	366	C26H54	000630-01-3	94
2		Eicosane	282	C20H42	000112-95-8	94
3		Nonadecane	268	C19H40	000629-92-5	93
4		Nonadecane	268	C19H40	000629-92-5	93
5		Octacosane, 2-methyl-	408	C29H60	001560-98-1	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN122022\
Data File : BN023296.D
Acq On : 20 Dec 2022 13:38
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ClientSampleId :
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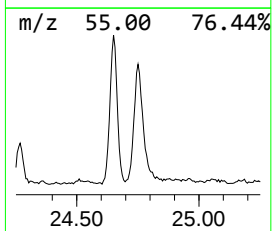
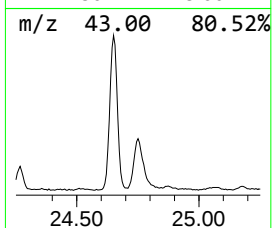
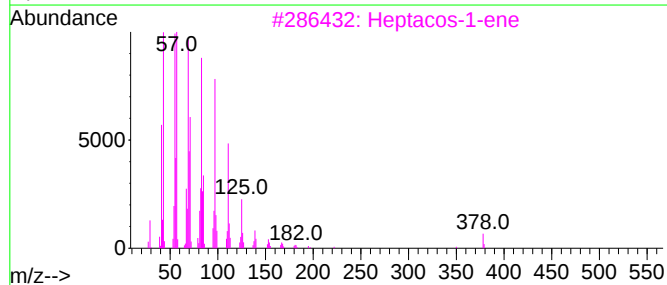
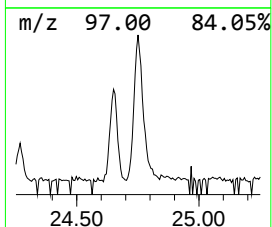
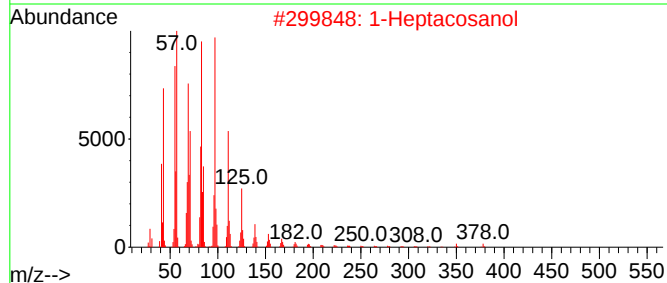
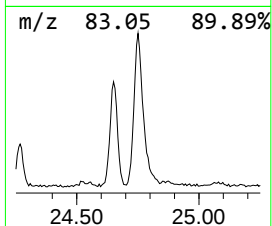
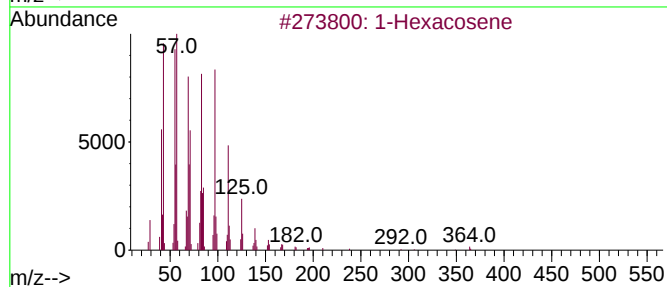
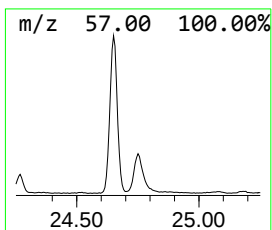
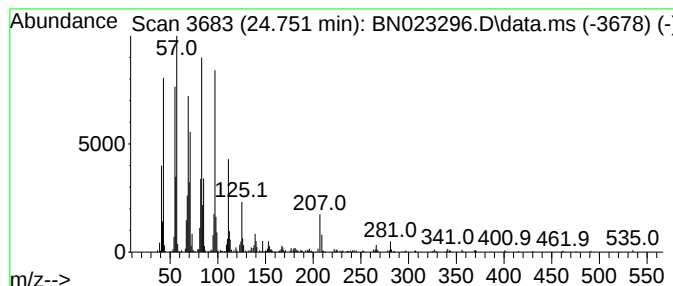
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.751	3.80 ng/ul	854579	Perylene-d12	24.027

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Hexacosene			364	C26H52	018835-33-1	95
2	1-Heptacosanol			396	C27H56O	002004-39-9	94
3	Heptacos-1-ene			378	C27H54	015306-27-1	94
4	Tricosyl trifluoroacetate			436	C25H47F3O2	1000351-75-1	93
5	1-Heneicosanol			312	C21H44O	015594-90-8	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN122022\
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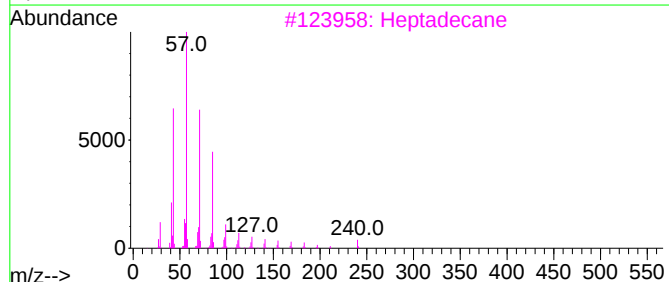
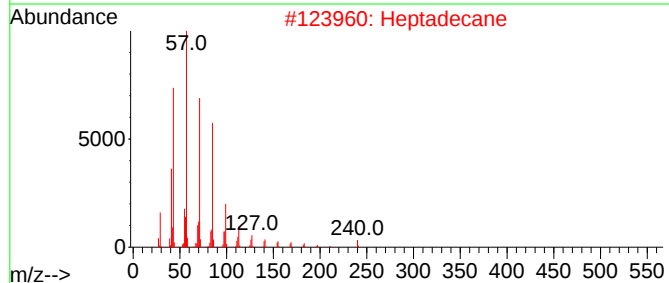
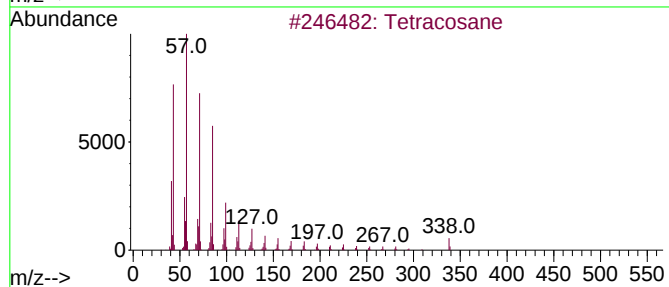
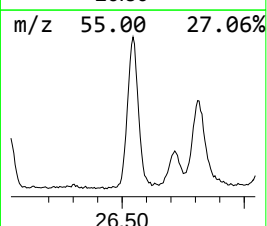
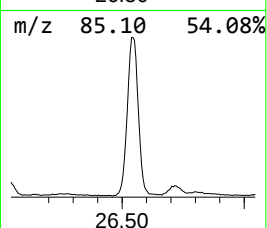
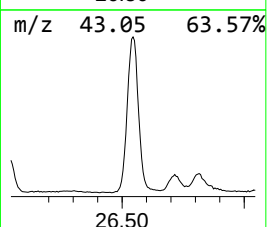
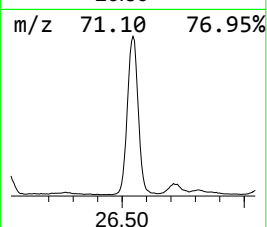
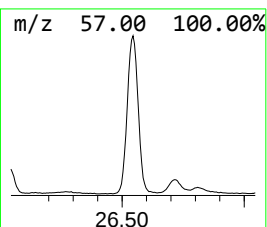
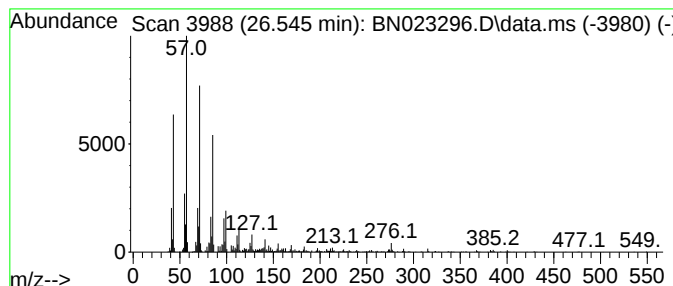
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.545	14.50 ng/ul	3259090	Perylene-d12	24.027

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetracosane	338	C24H50	000646-31-1	93
2			Heptadecane	240	C17H36	000629-78-7	93
3			Heptadecane	240	C17H36	000629-78-7	91
4			Heneicosane	296	C21H44	000629-94-7	90
5			Hexacosane, 1-iodo-	492	C26H53I	1000406-32-1	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN122022\
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Misc :
ALS Vial : 8 Sample Multiplier: 1

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ClientSampleId :
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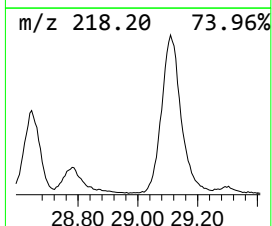
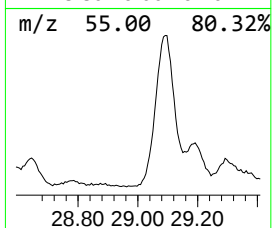
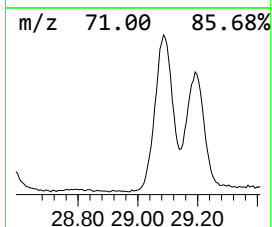
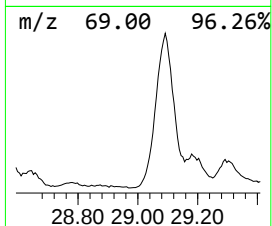
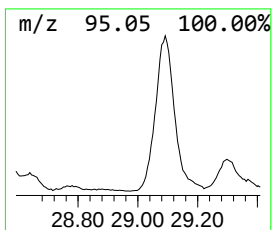
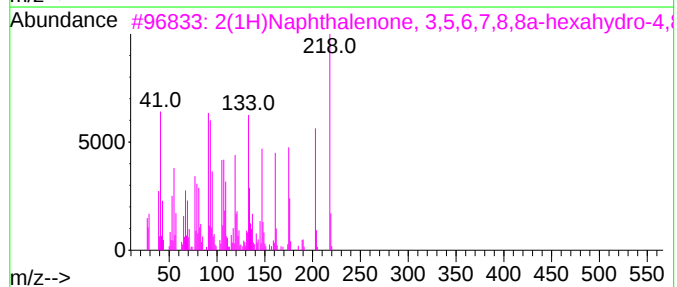
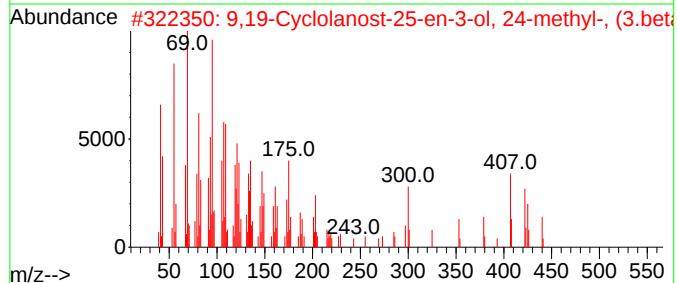
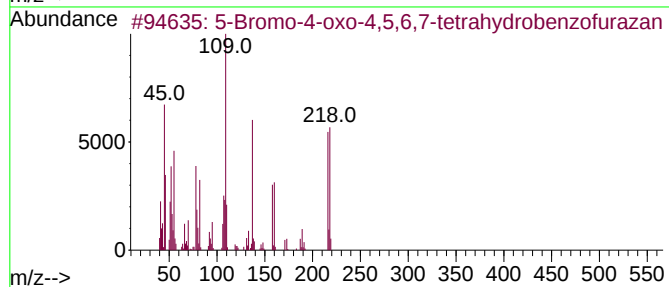
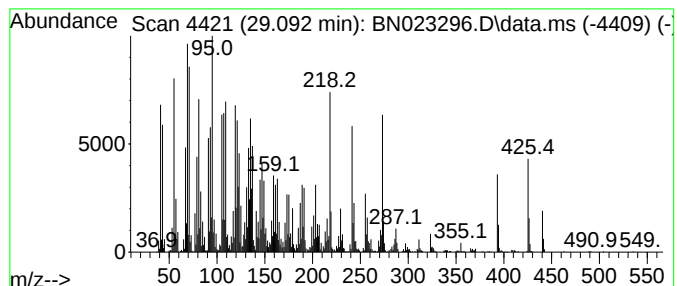
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN121522.M
Quant Title : SVOA CALIBRATION

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

Peak Number 11 5-Bromo-4-oxo-4,5,6,7-tetra... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
29.092	33.43 ng/ul	7514080	Perylene-d12	24.027

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			5-Bromo-4-oxo-4,5,6,7-tetrahydro...	216	C6H5BrN2O2	300574-36-1	93
2			9,19-Cyclolanost-25-en-3-ol, 24-...	440	C31H52O	000511-61-5	83
3			2(1H)Naphthalenone, 3,5,6,7,8,8a...	218	C15H22O	1000188-66-5	55
4			D:C-Friedo-B':A'-neogammacer-9(1...	440	C31H52O	004555-56-0	55
5			9,19-Cyclolanost-24-en-3-ol, (3...	426	C30H50O	000469-38-5	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN122022\
Data File : BN023296.D
Acq On : 20 Dec 2022 13:38
Operator : CG/JU
Sample : N6003-18
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
DBXT8

Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN121522.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
2-Pentanone, 4-...	5.064	4.7	ng/ul	388654	1	7.993	1669280	20.0
Benzophenone	16.004	12.7	ng/ul	2361120	3	14.645	3709520	20.0
n-Hexadecanoic ...	18.286	7.5	ng/ul	1699810	4	17.392	4548000	20.0
Benzene, hexaet...	18.869	2.1	ng/ul	486895	4	17.392	4548000	20.0
Octadecanoic acid	19.563	2.4	ng/ul	687426	5	21.580	5717860	20.0
4-(3-Fluoro-8-n...	19.945	12.0	ng/ul	3435390	5	21.580	5717860	20.0
1-Heneicosyl fo...	21.275	16.0	ng/ul	4586350	5	21.580	5717860	20.0
(DEL) Alkane: S...	24.651	6.7	ng/ul	1494160	6	24.027	4495080	20.0
(DEL) Alkane: S...	24.751	3.8	ng/ul	854579	6	24.027	4495080	20.0
(DEL) Alkane: S...	26.545	14.5	ng/ul	3259090	6	24.027	4495080	20.0
5-Bromo-4-oxo-4...	29.092	33.4	ng/ul	7514080	6	24.027	4495080	20.0