

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN122022\
 Data File : BN023295.D
 Acq On : 20 Dec 2022 13:01
 Operator : CG/JU
 Sample : N6003-17
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 DBXT1

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: BN023295.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	40	43	52	rVB	55544	83483	1.03%	0.075%
2	3.640	92	94	102	rVB	7813	10041	0.12%	0.009%
3	3.775	114	117	133	rBV	409121	683468	8.41%	0.614%
4	3.887	133	136	143	rVB2	20828	29826	0.37%	0.027%
5	4.017	153	158	165	rVB	19991	33297	0.41%	0.030%
6	4.111	171	174	180	rVB	17507	26613	0.33%	0.024%
7	5.064	330	336	347	rVB2	127706	197947	2.44%	0.178%
8	6.969	657	660	664	rBV5	8179	14296	0.18%	0.013%
9	7.152	683	691	701	rBV	1176767	1817414	22.36%	1.633%
10	7.322	712	720	730	rBV	957914	1439373	17.71%	1.293%
11	7.522	744	754	763	rBV	1187866	1844562	22.70%	1.657%
12	7.605	765	768	775	rVB6	8269	13508	0.17%	0.012%
13	7.993	825	834	842	rBV	1040070	1657490	20.39%	1.489%
14	8.058	842	845	852	rVB2	14590	25575	0.31%	0.023%
15	8.128	852	857	862	rBV	15653	25842	0.32%	0.023%
16	8.710	946	956	966	rBV	1471512	2380316	29.29%	2.138%
17	8.828	971	976	981	rVB	21761	33955	0.42%	0.031%
18	9.163	1024	1033	1041	rBV	1147173	1864267	22.94%	1.675%
19	9.269	1045	1051	1055	rVV8	6904	14006	0.17%	0.013%
20	9.334	1057	1062	1068	rVB9	5548	10696	0.13%	0.010%
21	9.581	1095	1104	1112	rVB2	20566	40273	0.50%	0.036%
22	9.893	1148	1157	1169	rBV	992597	1597621	19.66%	1.435%
23	10.110	1190	1194	1200	rVB8	7246	12251	0.15%	0.011%
24	10.434	1241	1249	1261	rBV	1535576	2572099	31.65%	2.311%
25	10.593	1272	1276	1280	rVB6	10981	17357	0.21%	0.016%
26	10.816	1306	1314	1326	rVB	1607582	2649577	32.60%	2.380%
27	10.952	1330	1337	1353	rBV	1425019	2369085	29.15%	2.128%
28	11.204	1375	1380	1389	rVB2	21273	41767	0.51%	0.038%
29	11.293	1389	1395	1399	rBV2	13781	22044	0.27%	0.020%
30	11.604	1441	1448	1454	rBV7	9796	20251	0.25%	0.018%
31	12.234	1549	1555	1560	rBV2	16831	26485	0.33%	0.024%
32	12.399	1578	1583	1590	rVB	30459	45966	0.57%	0.041%
33	13.387	1746	1751	1755	rBV7	6469	12016	0.15%	0.011%
34	13.469	1760	1765	1772	rVV	32422	48655	0.60%	0.044%
35	13.540	1772	1777	1786	rVV7	8603	20456	0.25%	0.018%
36	13.640	1786	1794	1801	rVB	97035	134323	1.65%	0.121%

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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-BN121522.M
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37	13.969	1846	1850	1852	rBV3	8222	10662	0.13%	0.010%
38	14.057	1858	1865	1874	rBV	2859348	3994906	49.15%	3.589%
39	14.122	1874	1876	1880	rVV4	9107	10839	0.13%	0.010%
40	14.175	1880	1885	1889	rVB6	13080	21883	0.27%	0.020%
41	14.240	1891	1896	1900	rBV2	11476	17054	0.21%	0.015%
42	14.340	1906	1913	1923	rVV	3433048	4898837	60.27%	4.401%
43	14.534	1942	1946	1951	rVB2	21093	28709	0.35%	0.026%
44	14.645	1957	1965	1973	rBV	2596823	3714965	45.71%	3.338%
45	14.840	1990	1998	2013	rBV	1363295	2027024	24.94%	1.821%
46	14.992	2021	2024	2029	rVB5	11463	18478	0.23%	0.017%
47	15.057	2030	2035	2040	rVB5	37498	57251	0.70%	0.051%
48	15.428	2093	2098	2102	rBV6	6612	13481	0.17%	0.012%
49	15.487	2105	2108	2115	rVB4	9696	14212	0.17%	0.013%
50	15.640	2125	2134	2141	rBV	4270513	6175687	75.98%	5.548%
51	15.751	2146	2153	2165	rVV	1393020	1853043	22.80%	1.665%
52	15.875	2169	2174	2183	rVB2	22916	38021	0.47%	0.034%
53	16.004	2186	2196	2201	rBV	1546432	2122181	26.11%	1.907%
54	16.351	2252	2255	2262	rVB7	10015	17850	0.22%	0.016%
55	16.416	2264	2266	2274	rVB9	7141	12341	0.15%	0.011%
56	16.516	2276	2283	2287	rBV8	6450	15680	0.19%	0.014%
57	16.563	2287	2291	2297	rVB	49722	63762	0.78%	0.057%
58	16.787	2325	2329	2337	rVB7	9790	21905	0.27%	0.020%
59	17.181	2392	2396	2402	rVB4	7861	16886	0.21%	0.015%
60	17.281	2410	2413	2415	rBV4	9875	12159	0.15%	0.011%
61	17.392	2425	2432	2442	rBV	3249669	4592651	56.51%	4.126%
62	17.492	2442	2449	2458	rVV	4604670	6639346	81.69%	5.965%
63	17.575	2459	2463	2469	rVB6	36636	56526	0.70%	0.051%
64	17.669	2475	2479	2487	rVB2	16769	33973	0.42%	0.031%
65	17.769	2491	2496	2503	rBV5	20926	43581	0.54%	0.039%
66	18.057	2542	2545	2550	rVB2	24314	29572	0.36%	0.027%
67	18.163	2558	2563	2569	rBV8	17671	36997	0.46%	0.033%
68	18.228	2570	2574	2579	rBV	35901	60253	0.74%	0.054%
69	18.286	2579	2584	2593	rBV	611471	916115	11.27%	0.823%
70	18.481	2610	2617	2622	rBV	130548	176216	2.17%	0.158%
71	18.533	2622	2626	2631	rVB	98761	128173	1.58%	0.115%
72	18.710	2651	2656	2663	rBV3	260892	319580	3.93%	0.287%
73	18.798	2667	2671	2677	rVV	97996	125500	1.54%	0.113%
74	18.863	2678	2682	2690	rVV	288496	388710	4.78%	0.349%
75	19.145	2726	2730	2733	rBV5	18751	29619	0.36%	0.027%
76	19.216	2737	2742	2746	rVV	272062	330808	4.07%	0.297%

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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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77	19.351	2761	2765	2769	rBV4	79001	99137	1.22%	0.089%
78	19.416	2772	2776	2778	rBV	118014	141049	1.74%	0.127%
79	19.445	2778	2781	2784	rVB2	72199	87899	1.08%	0.079%
80	19.563	2796	2801	2807	rBV	250842	401000	4.93%	0.360%
81	19.704	2821	2825	2830	rVB5	47386	69220	0.85%	0.062%
82	19.786	2833	2839	2849	rVV	6053792	8127597	100.00%	7.302%
83	19.945	2861	2866	2874	rBV	3306762	3833188	47.16%	3.444%
84	20.069	2883	2887	2893	rBV	137483	202222	2.49%	0.182%
85	20.128	2894	2897	2905	rVB2	140225	190500	2.34%	0.171%
86	20.298	2921	2926	2933	rBV2	507177	693985	8.54%	0.623%
87	20.539	2964	2967	2975	rVB8	68197	101953	1.25%	0.092%
88	20.816	3011	3014	3018	rVB2	104693	117952	1.45%	0.106%
89	21.280	3075	3093	3112	rBV2	1662891	5640852	69.40%	5.068%
90	21.580	3138	3144	3148	rBV	4161950	6220244	76.53%	5.588%
91	21.727	3167	3169	3175	rVB3	160298	202179	2.49%	0.182%
92	22.198	3246	3249	3252	rBV2	134908	206970	2.55%	0.186%
93	22.704	3332	3335	3341	rVB2	167019	220097	2.71%	0.198%
94	23.269	3426	3431	3436	rBV	454804	721670	8.88%	0.648%
95	23.868	3526	3533	3548	rVB	3400512	7544170	92.82%	6.778%
96	24.027	3553	3560	3577	rVB	2203636	4490594	55.25%	4.034%
97	24.651	3660	3666	3673	rVB	593154	1214711	14.95%	1.091%
98	24.751	3679	3683	3696	rVB2	299992	710275	8.74%	0.638%
99	26.539	3980	3987	4000	rVB	838988	2559872	31.50%	2.300%
100	29.092	4410	4421	4434	rBV6	1625082	6590690	81.09%	5.921%

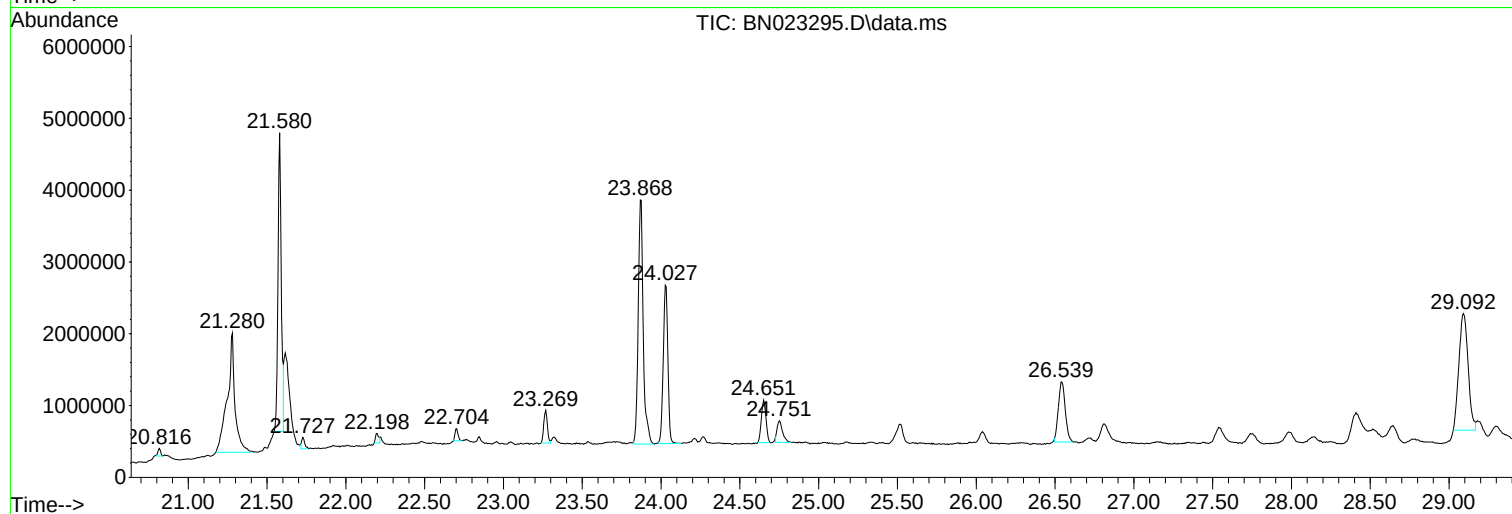
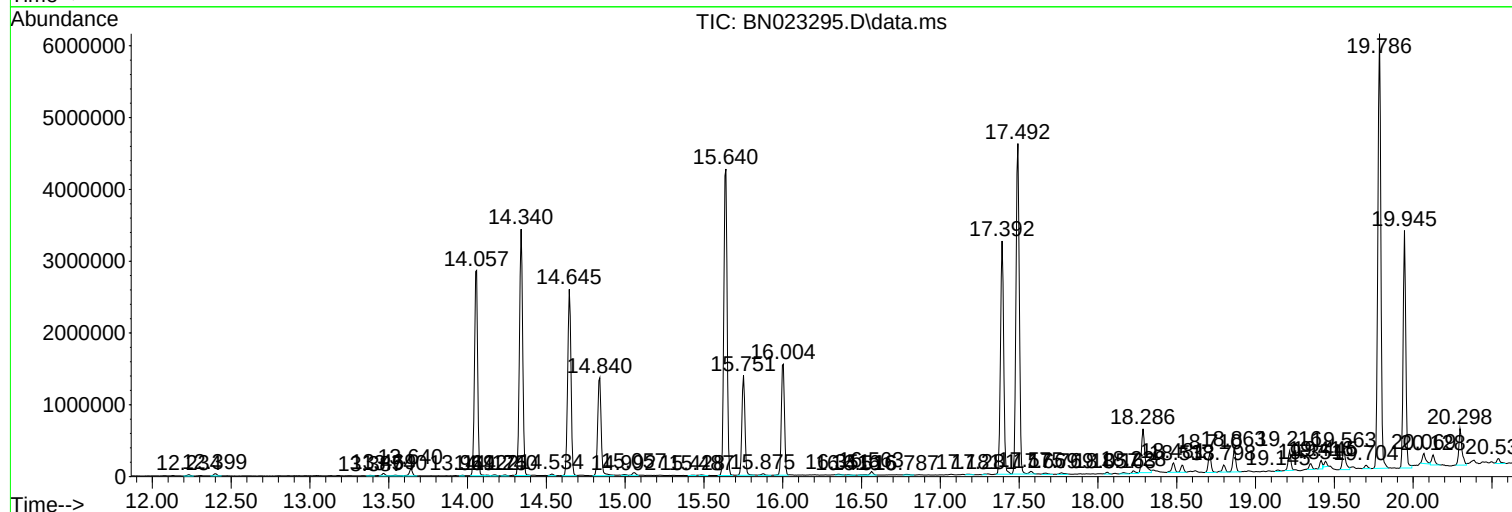
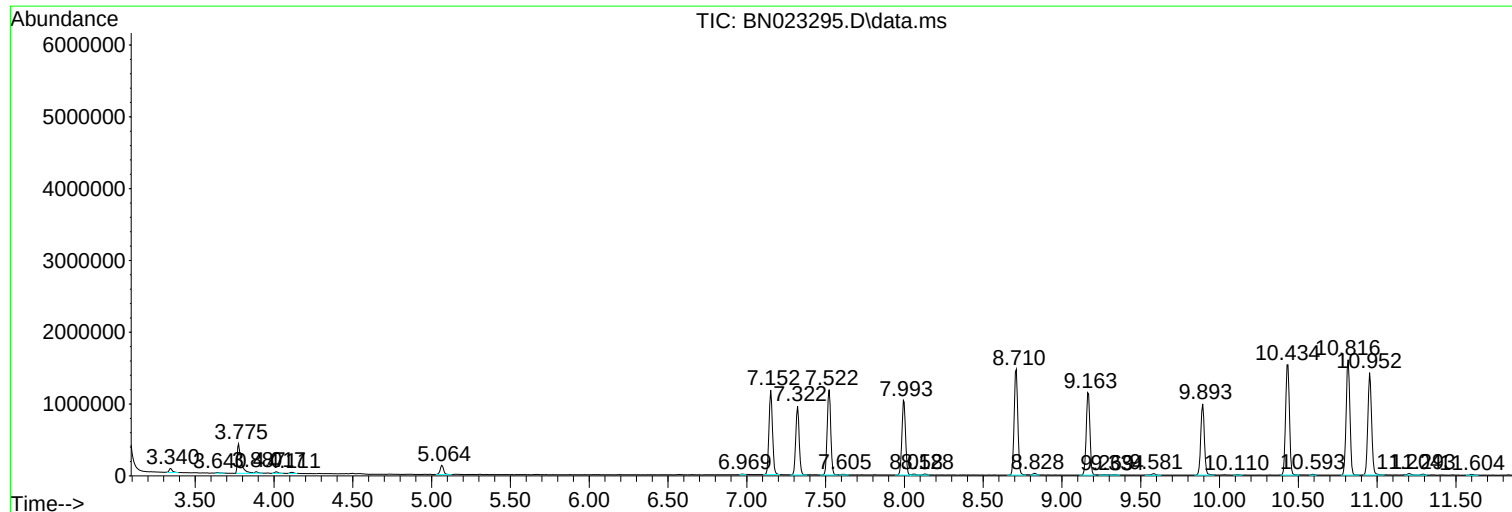
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Instrument :
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DBXT1

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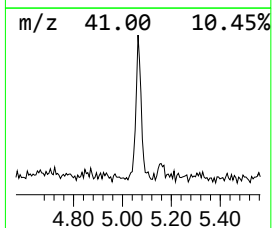
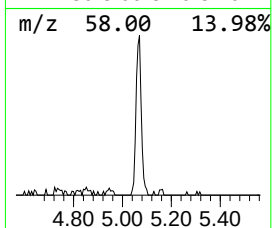
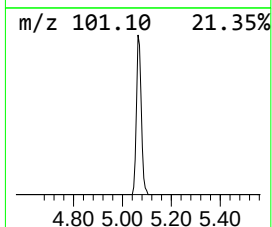
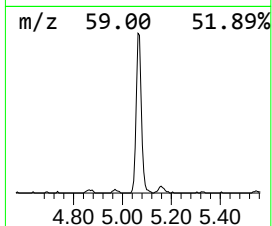
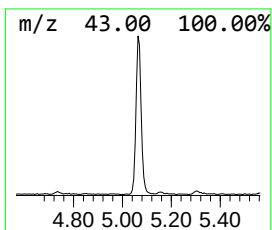
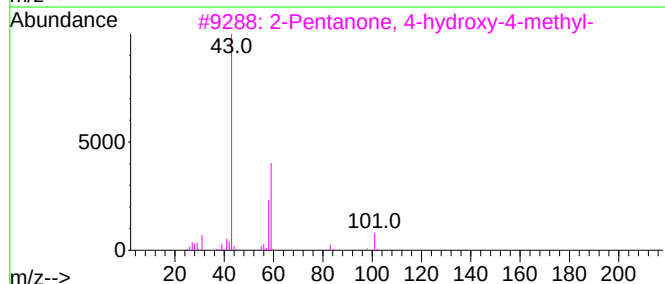
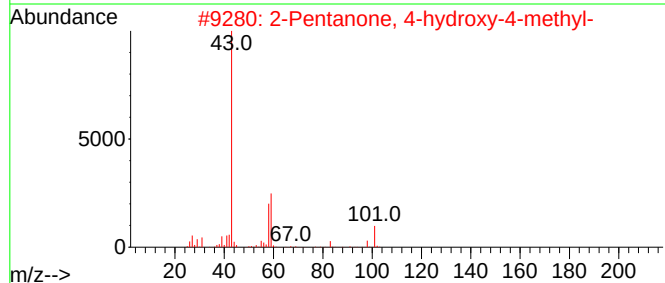
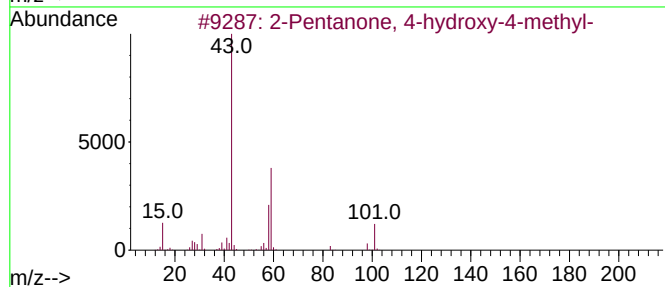
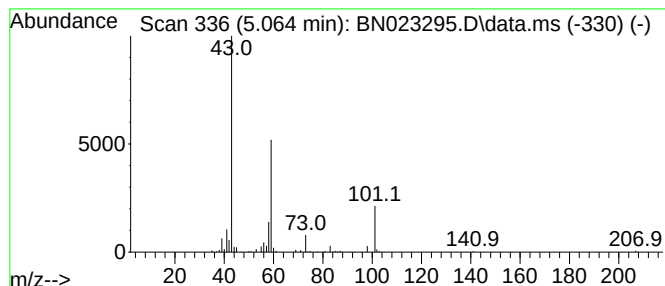
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN121522.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.064	2.39 ng/ul	197947	1,4-Dichlorobenzene-d4	7.999

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50		
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	50		
4	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	38		
5	Propane, 1-ethoxy-2-methyl-	102	C6H14O	000627-02-1	25		



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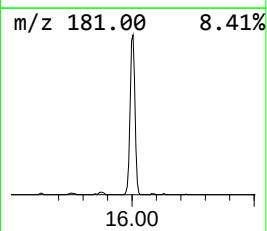
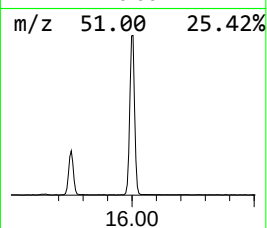
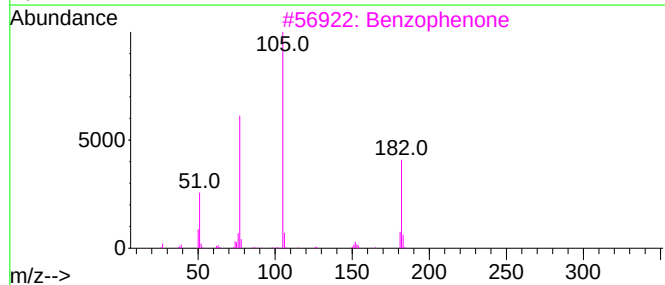
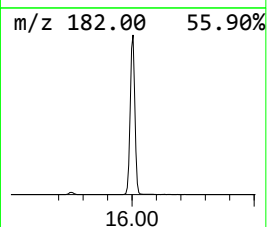
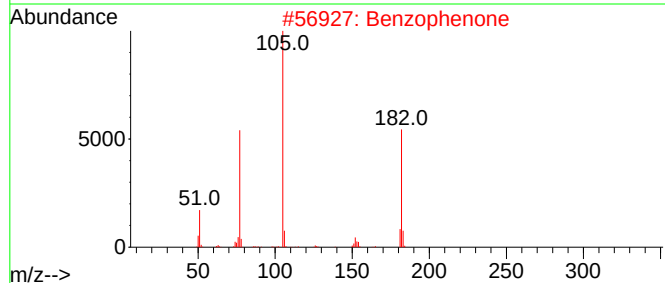
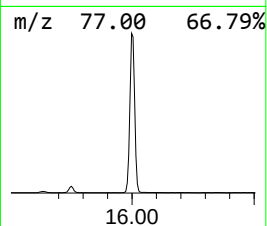
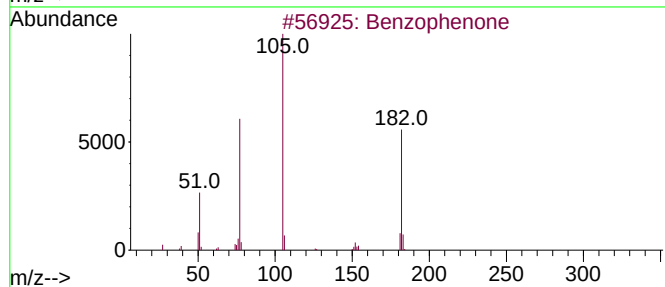
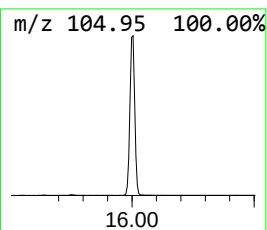
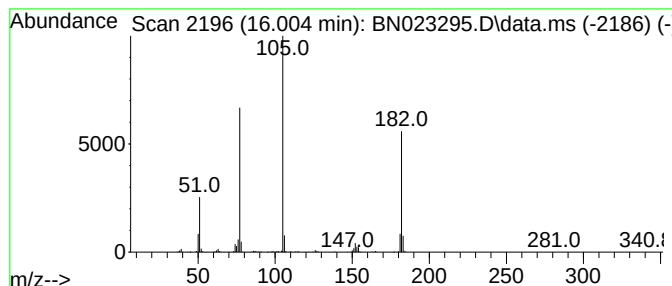
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN121522.M
Quant Title : SVOA CALIBRATION

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	11.43 ng/ul	2122180	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	91
5			Benzophenone	182	C13H10O	000119-61-9	90



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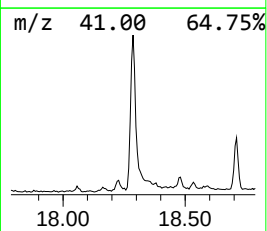
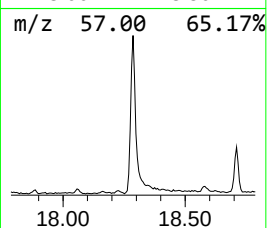
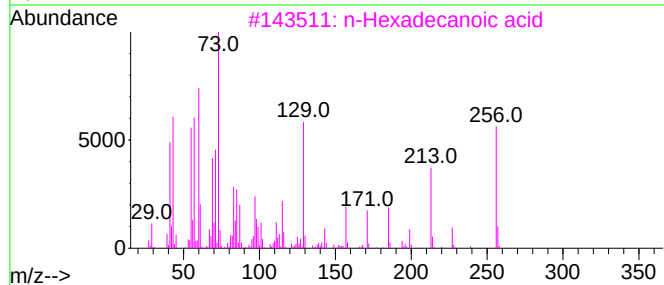
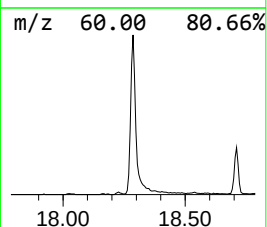
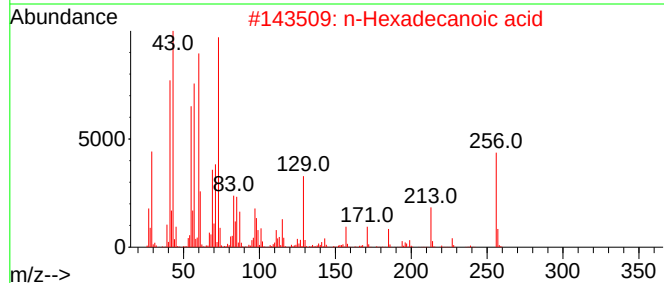
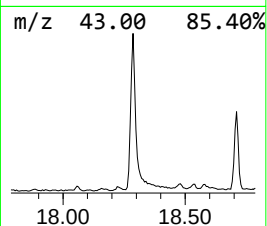
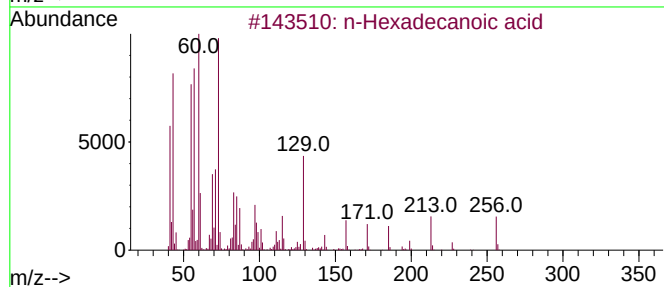
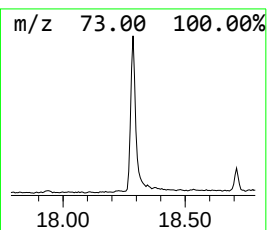
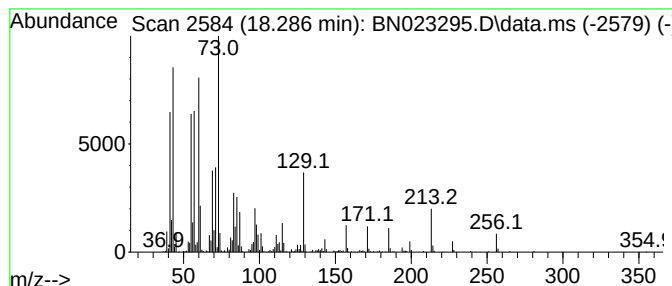
Instrument :
BNA_N
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN121522.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 n-Hexadecanoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.	
18.286	3.99 ng/ul	916115	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	98
3	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
4	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	97
5	Octadecanoic acid	284	C18H36O2	000057-11-4	93



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Data File : BN023295.D
Acq On : 20 Dec 2022 13:01
Operator : CG/JU
Sample : N6003-17
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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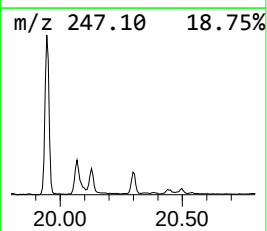
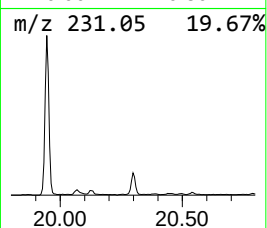
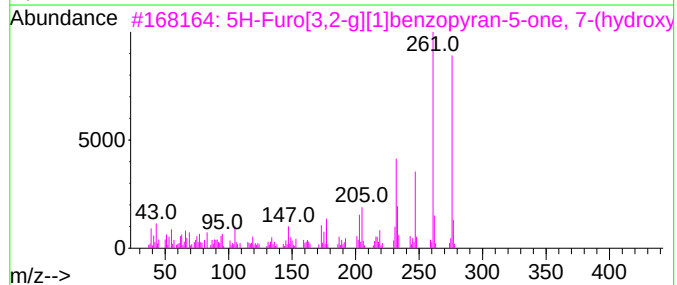
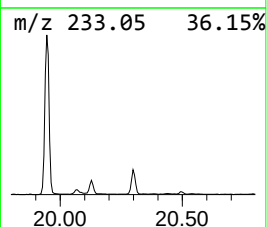
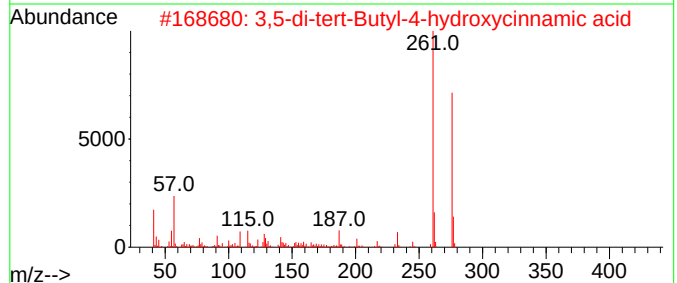
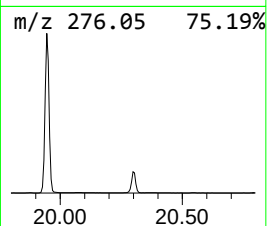
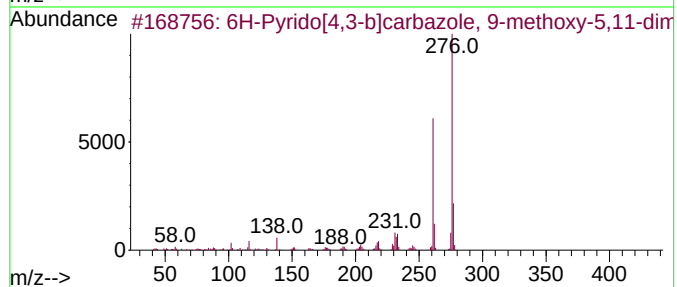
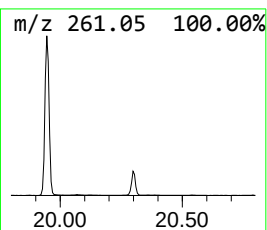
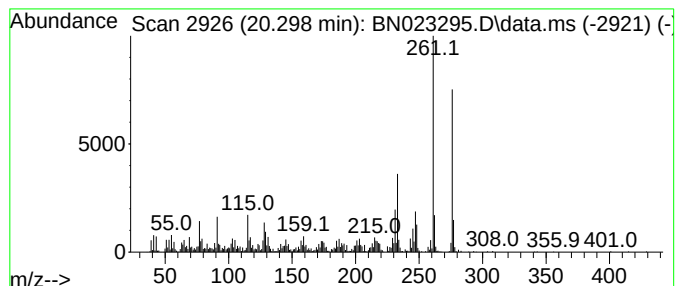
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 6H-Pyrido[4,3-b]carbazole, ... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.298	2.23 ng/ul	693985	Chrysene-d12	21.580

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			6H-Pyrido[4,3-b]carbazole, 9-met...	276	C18H16N2O	010371-86-5	83
2			3,5-di-tert-Butyl-4-hydroxycinna...	276	C17H24O3	022014-01-3	76
3			5H-Furo[3,2-g][1]benzopyran-5-on...	276	C14H12O6	000668-10-0	72
4			3-Methoxy-5-pentyl-2-prenylpheno...	276	C18H28O2	083036-53-7	68
5			3-(4-Fluorophenoxy)phenol, TMS	276	C15H17F02Si	1000495-60-9	58



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Sample : N6003-17
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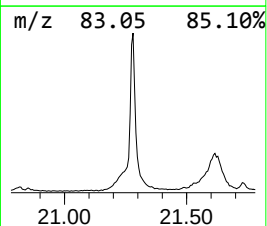
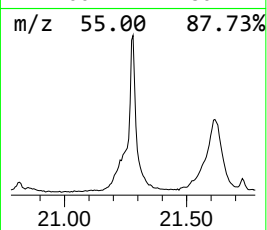
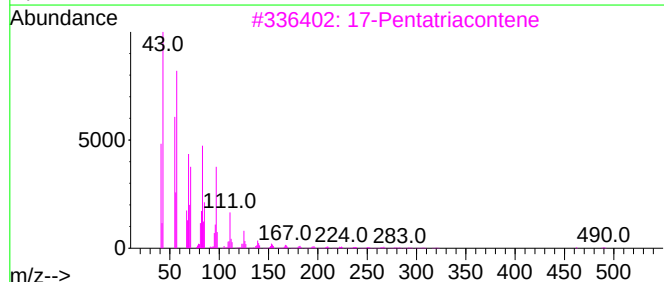
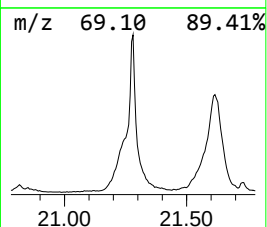
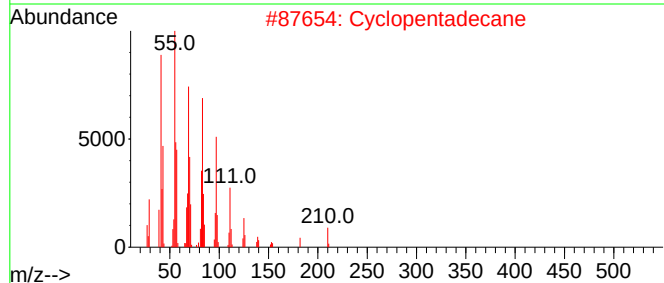
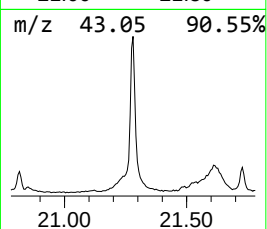
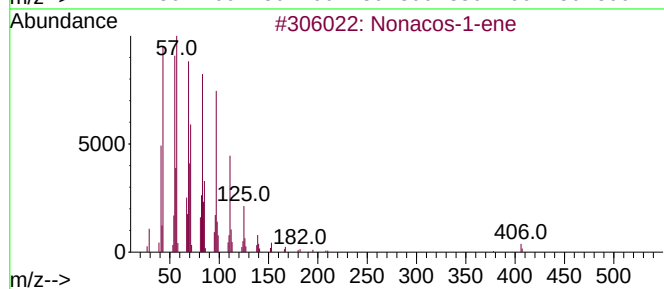
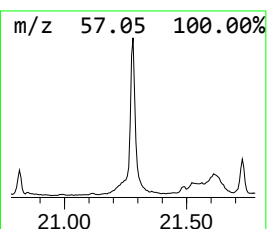
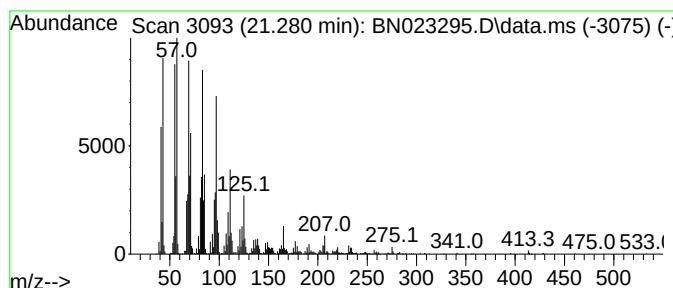
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD		R.T.	
21.280	18.14 ng/ul	5640850	Chrysene-d12		21.580	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Nonacos-1-ene		406	C29H58	018835-35-3	91
2	Cyclopentadecane		210	C15H30	000295-48-7	90
3	17-Pentatriacontene		491	C35H70	006971-40-0	90
4	Heptacos-1-ene		378	C27H54	015306-27-1	87
5	Triacontan-15-ol		438	C30H62O	031849-26-0	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN122022\
Data File : BN023295.D
Acq On : 20 Dec 2022 13:01
Operator : CG/JU
Sample : N6003-17
Misc :
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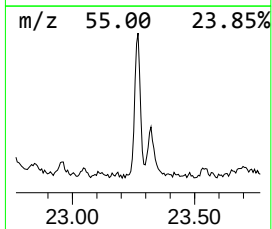
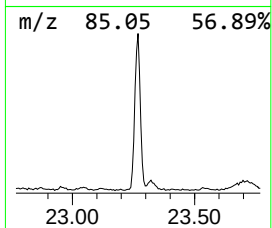
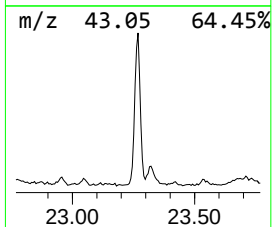
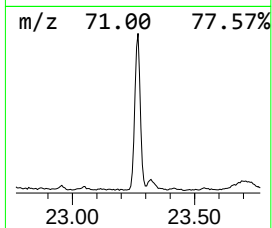
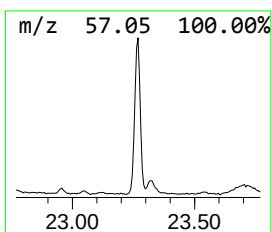
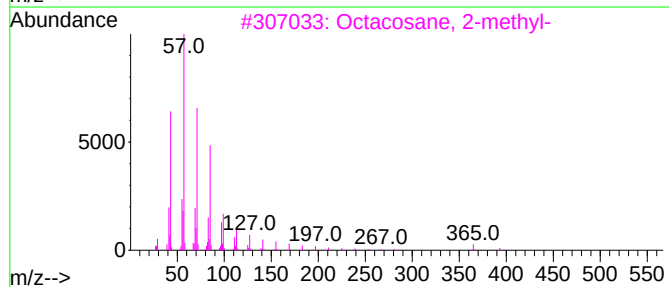
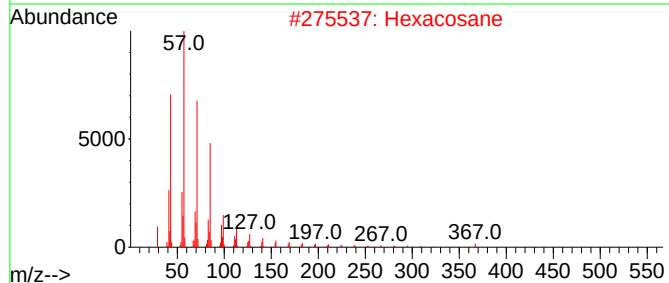
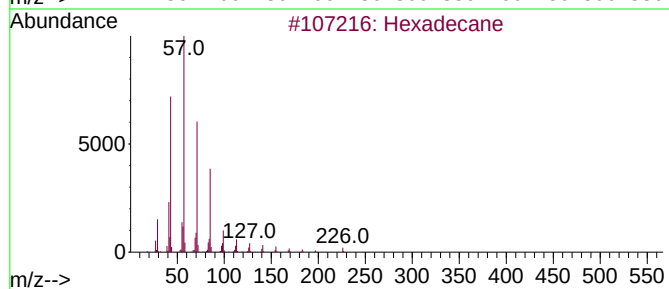
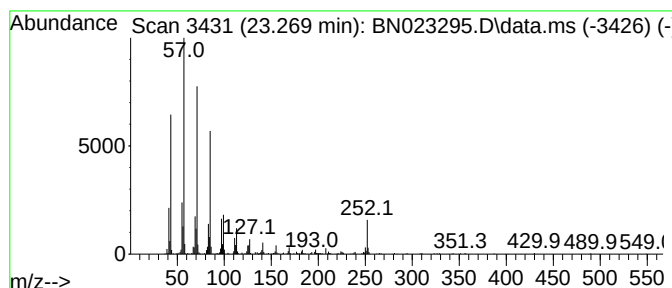
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.269	3.21 ng/ul	721670	Perylene-d12	24.033

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexadecane	226	C16H34	000544-76-3	95
2			Hexacosane	366	C26H54	000630-01-3	93
3			Octacosane, 2-methyl-	408	C29H60	001560-98-1	90
4			Tetracosane	338	C24H50	000646-31-1	90
5			Octacosane	394	C28H58	000630-02-4	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN122022\
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Sample : N6003-17
Misc :
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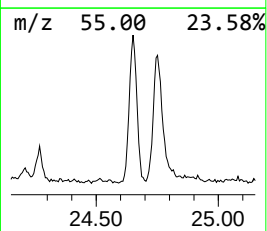
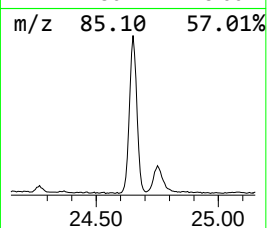
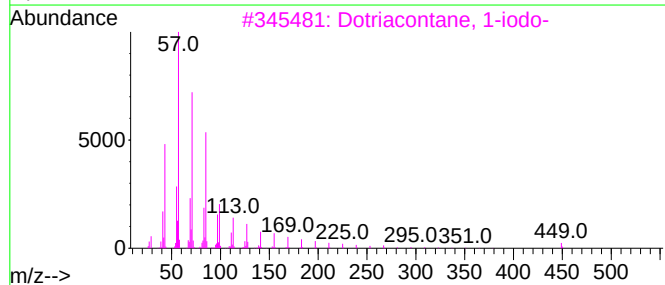
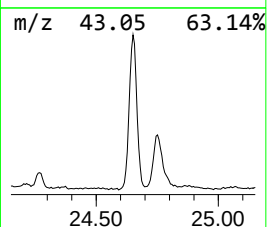
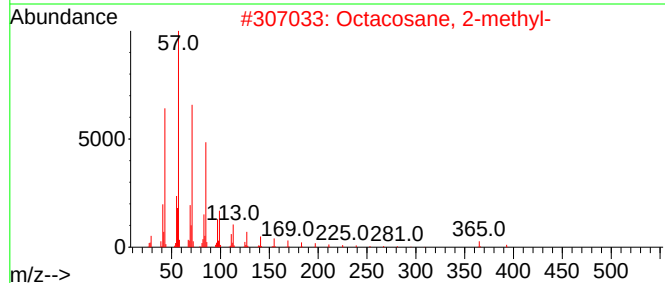
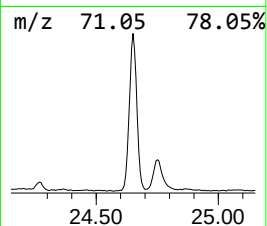
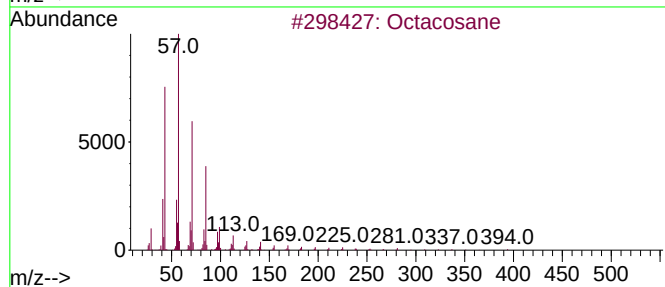
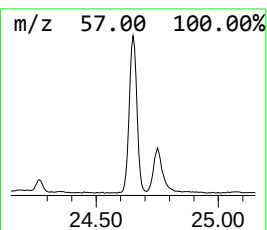
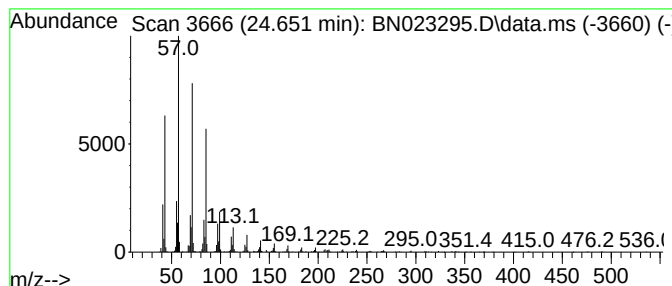
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
24.651	5.41 ng/ul	1214710	Perylene-d12	24.033	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octacosane	394	C28H58	000630-02-4	91
2	Octacosane, 2-methyl-	408	C29H60	001560-98-1	91
3	Dotriacontane, 1-iodo-	576	C32H65I	1000406-32-4	91
4	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
5	Heneicosane	296	C21H44	000629-94-7	91



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Data File : BN023295.D
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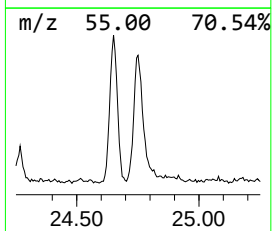
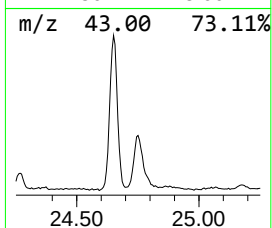
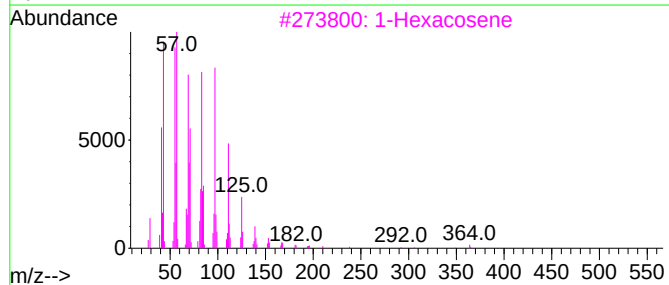
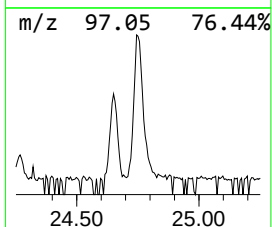
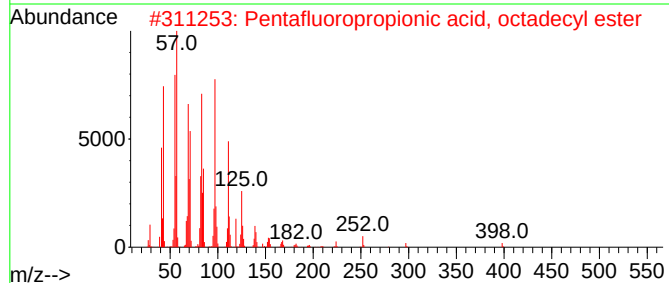
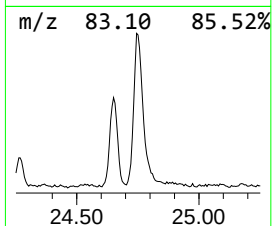
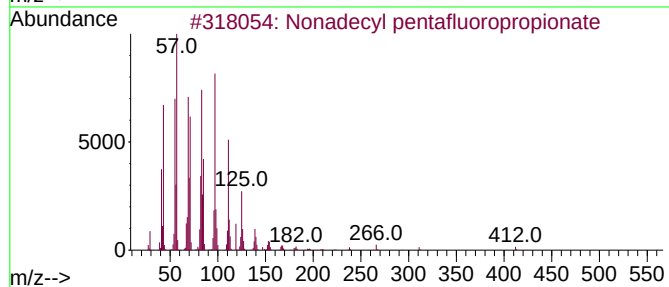
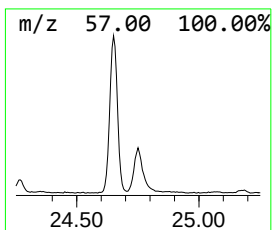
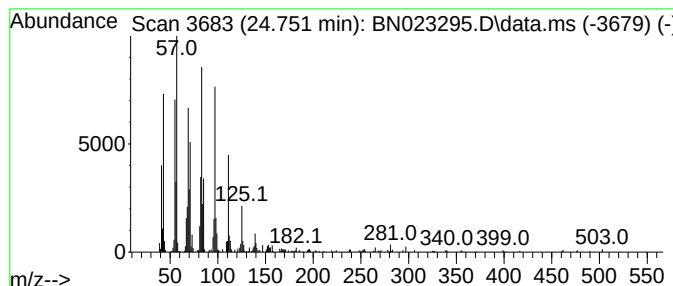
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 Nonadecyl pentafluoropropio... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.751	3.16 ng/ul	710275	Perylene-d12	24.033

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonadecyl pentafluoropropionate	430	C22H39F5O2	1000351-88-8	94
2			Pentafluoropropionic acid, octadecyl ester	416	C21H37F5O2	959261-25-7	94
3			1-Hexacosene	364	C26H52	018835-33-1	93
4			Heptafluorobutyric acid, n-octadecyl ester	466	C22H37F7O2	000400-57-7	91
5			Octacosanol	410	C28H58O	000557-61-9	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN122022\
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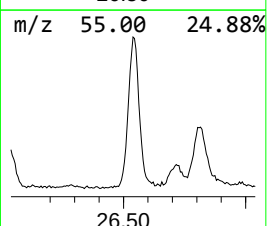
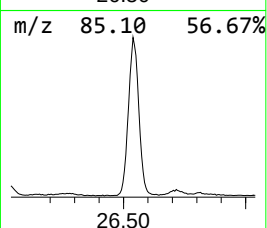
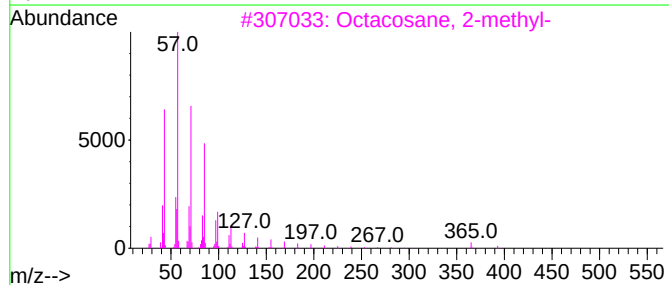
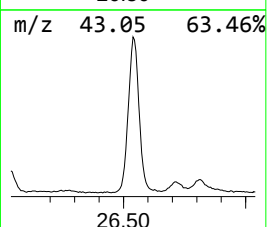
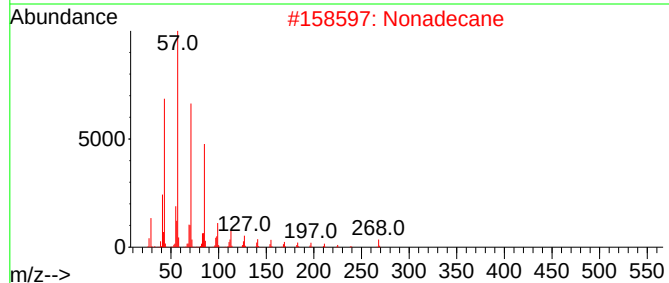
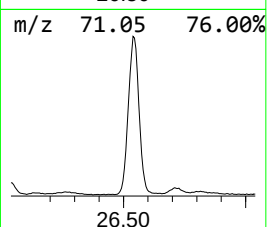
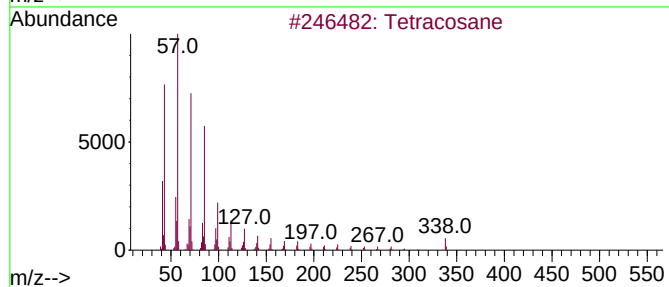
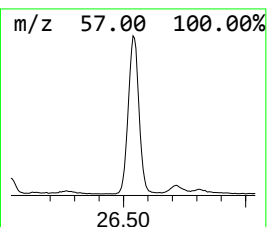
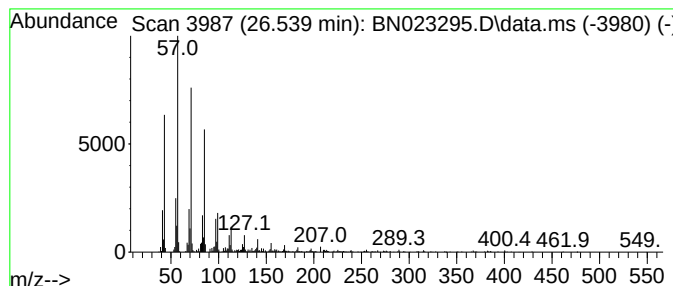
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD		R.T.	
26.539	11.40 ng/ul	2559870	Perylene-d12		24.033	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Tetracosane		338	C24H50	000646-31-1	97
2	Nonadecane		268	C19H40	000629-92-5	93
3	Octacosane, 2-methyl-		408	C29H60	001560-98-1	91
4	Heptacosane		380	C27H56	000593-49-7	91
5	2-Methyltetracosane		352	C25H52	001560-78-7	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN122022\
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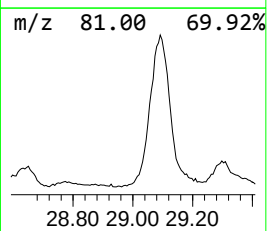
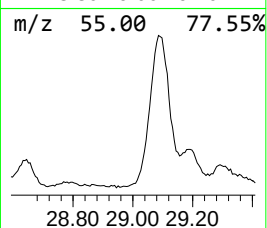
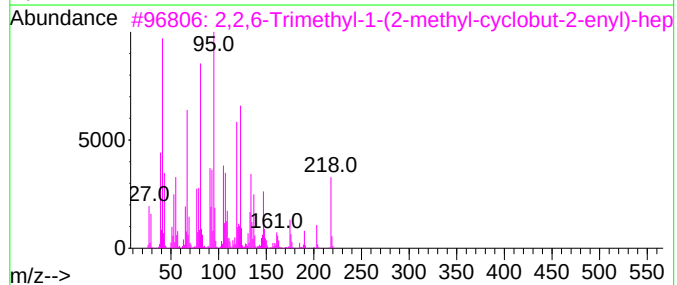
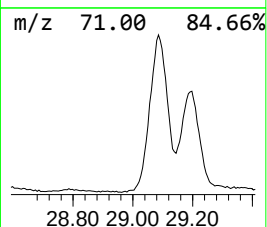
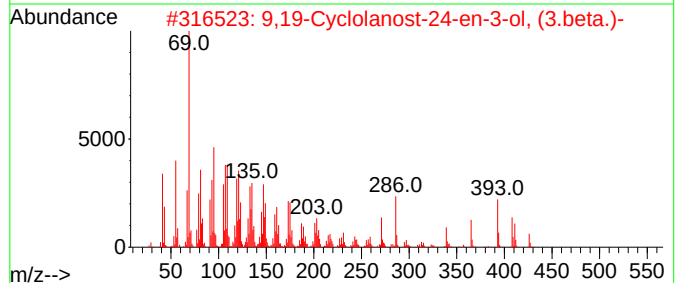
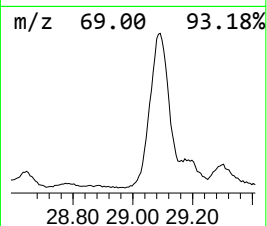
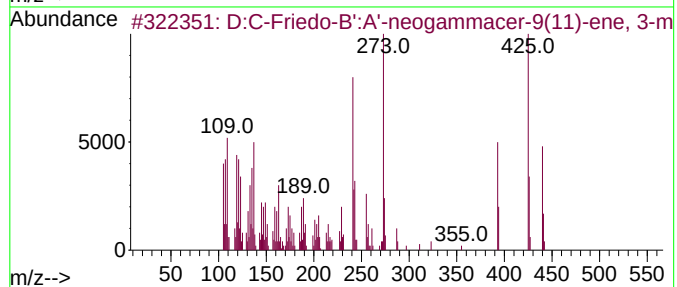
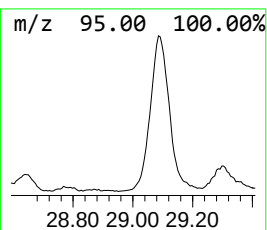
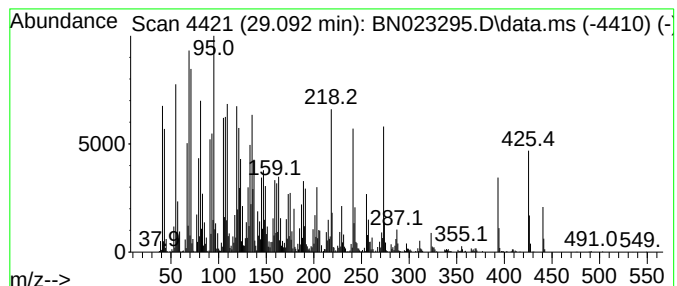
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BNA_N
ClientSampleId :
DBXT1

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 D:C-Friedo-B':A'-neogammace... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.		
29.092	29.35 ng/ul	6590690	Perylene-d12	24.033		
Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		D:C-Friedo-B':A'-neogammacer-9(1...	440	C31H52O	004555-56-0	93
2		9,19-Cyclolanost-24-en-3-ol, (3...	426	C30H50O	000469-38-5	53
3		2,2,6-Trimethyl-1-(2-methyl-cycl...	218	C15H22O	1000188-72-8	35
4		9-Thiabicyclo[3.3.1]nonane, 2,6-...	274	C14H18N4S	1000141-52-1	35
5		6-Isopropenyl-4,8a-dimethyl-4a,5...	218	C15H22O	086917-79-5	25



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN122022\
 Data File : BN023295.D
 Acq On : 20 Dec 2022 13:01
 Operator : CG/JU
 Sample : N6003-17
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 DBXT1

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	2.4	ng/ul	197947	1	7.999	1657490	20.0
Benzophenone	16.004	11.4	ng/ul	2122180	3	14.645	3714970	20.0
n-Hexadecanoic ...	18.286	4.0	ng/ul	916115	4	17.392	4592650	20.0
4-(3-Fluoro-8-n...	19.945	12.3	ng/ul	3833190	5	21.580	6220240	20.0
6H-Pyrido[4,3-b...	20.298	2.2	ng/ul	693985	5	21.580	6220240	20.0
(DEL) Alkane: S...	21.280	18.1	ng/ul	5640850	5	21.580	6220240	20.0
(DEL) Alkane: S...	23.269	3.2	ng/ul	721670	6	24.033	4490590	20.0
(DEL) Alkane: S...	24.651	5.4	ng/ul	1214710	6	24.033	4490590	20.0
Nonadecyl penta...	24.751	3.2	ng/ul	710275	6	24.033	4490590	20.0
(DEL) Alkane: S...	26.539	11.4	ng/ul	2559870	6	24.033	4490590	20.0
D:C-Friedo-B':A...	29.092	29.4	ng/ul	6590690	6	24.033	4490590	20.0