

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070223\
 Data File : BP015926.D
 Acq On : 02 Jul 2023 16:18
 Operator : MA/JU
 Sample : 03243-08
 Misc :
 ALS Vial : 91 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DCGC8

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_P\Methods\SFAM-EPA-BP062623.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP015926.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.375	27	32	37	rVB	75275	102629	1.87%	0.119%
2	3.828	106	109	122	rBV	290635	536468	9.78%	0.620%
3	3.922	122	125	135	rVB2	33695	62490	1.14%	0.072%
4	4.034	140	144	152	rVB	91074	147052	2.68%	0.170%
5	4.140	159	162	167	rVB	21574	30042	0.55%	0.035%
6	4.540	226	230	234	rVB7	7337	11564	0.21%	0.013%
7	4.963	298	302	305	rBV4	11634	16906	0.31%	0.020%
8	5.087	319	323	328	rBV	14139	23892	0.44%	0.028%
9	5.816	441	447	465	rBV	170535	475874	8.68%	0.550%
10	6.005	477	479	483	rVB4	6693	7928	0.14%	0.009%
11	6.687	587	595	602	rBV	101177	160595	2.93%	0.186%
12	6.999	643	648	652	rBV2	33961	50122	0.91%	0.058%
13	7.234	681	688	700	rBV	823873	2013706	36.71%	2.328%
14	7.369	706	711	722	rBV	988567	1775577	32.37%	2.053%
15	7.452	722	725	739	rVB	91514	187093	3.41%	0.216%
16	7.581	740	747	774	rVB	1217047	2277759	41.53%	2.633%
17	8.046	819	826	843	rBV	1285227	2343597	42.73%	2.709%
18	8.240	854	859	866	rBV3	33705	69072	1.26%	0.080%
19	8.310	866	871	873	rVV	35499	58948	1.07%	0.068%
20	8.340	873	876	883	rVV5	30282	49763	0.91%	0.058%
21	8.440	886	893	898	rVB5	8437	17547	0.32%	0.020%
22	8.552	909	912	916	rVB6	5121	8389	0.15%	0.010%
23	8.787	945	952	971	rBV	900216	2201085	40.13%	2.545%
24	9.240	1021	1029	1052	rBV	1057213	2258140	41.17%	2.611%
25	9.387	1052	1054	1061	rVB8	7300	12626	0.23%	0.015%
26	9.575	1082	1086	1091	rVB4	11705	19846	0.36%	0.023%
27	9.634	1091	1096	1099	rVB7	4522	7855	0.14%	0.009%
28	9.969	1146	1153	1178	rBV	729762	1791195	32.66%	2.071%
29	10.487	1233	1241	1242	rBV8	6431	10685	0.19%	0.012%
30	10.540	1242	1250	1279	rVV	748482	2486906	45.34%	2.875%
31	10.875	1299	1307	1324	rVV	1755378	3394037	61.88%	3.924%
32	11.010	1325	1330	1332	rVV2	26917	45167	0.82%	0.052%
33	11.057	1332	1338	1364	rVB	297878	1009550	18.41%	1.167%
34	11.334	1379	1385	1396	rBV	35062	114111	2.08%	0.132%
35	11.428	1397	1401	1411	rVV4	20618	48743	0.89%	0.056%
36	11.869	1470	1476	1478	rBV6	6307	11194	0.20%	0.013%

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
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37	12.075	1506	1511	1516	rVB7	8432	14594	0.27%	0.017%
38	12.216	1528	1535	1540	rBV2	35138	58072	1.06%	0.067%
39	12.275	1540	1545	1555	rVB2	7545	18880	0.34%	0.022%
40	12.451	1570	1575	1579	rBV7	23121	44269	0.81%	0.051%
41	12.493	1579	1582	1590	rVB2	18403	38040	0.69%	0.044%
42	12.751	1621	1626	1635	rBV6	43204	89703	1.64%	0.104%
43	13.045	1671	1676	1679	rBV7	6199	13210	0.24%	0.015%
44	13.210	1699	1704	1706	rBV6	6206	9006	0.16%	0.010%
45	13.445	1739	1744	1745	rVV5	6588	10159	0.19%	0.012%
46	13.492	1748	1752	1757	rVV	24345	39955	0.73%	0.046%
47	13.557	1757	1763	1773	rVV6	36403	101162	1.84%	0.117%
48	13.645	1773	1778	1779	rVV5	6280	11691	0.21%	0.014%
49	13.675	1779	1783	1790	rVB8	15749	27610	0.50%	0.032%
50	13.998	1833	1838	1844	rBV4	58286	94930	1.73%	0.110%
51	14.110	1850	1857	1869	rBV	2056628	3468570	63.24%	4.010%
52	14.263	1882	1883	1889	rVB6	7713	8338	0.15%	0.010%
53	14.392	1898	1905	1921	rBV2	2631473	4383833	79.93%	5.068%
54	14.510	1922	1925	1929	rVV6	21145	41548	0.76%	0.048%
55	14.563	1929	1934	1942	rVB2	50744	84610	1.54%	0.098%
56	14.634	1942	1946	1951	rBV3	33925	50819	0.93%	0.059%
57	14.698	1951	1957	1978	rVB	2514955	4069295	74.19%	4.705%
58	14.922	1989	1995	2001	rBV3	44705	65944	1.20%	0.076%
59	15.016	2004	2011	2023	rBV2	121863	490340	8.94%	0.567%
60	15.257	2048	2052	2057	rVB5	21194	33899	0.62%	0.039%
61	15.463	2083	2087	2091	rBV5	11549	18272	0.33%	0.021%
62	15.516	2093	2096	2101	rVB2	21220	28328	0.52%	0.033%
63	15.598	2103	2110	2116	rBV3	69495	107417	1.96%	0.124%
64	15.698	2119	2127	2146	rBV	2734326	4957174	90.38%	5.731%
65	15.875	2152	2157	2179	rBV	264925	806361	14.70%	0.932%
66	16.069	2185	2190	2198	rBV	330462	558475	10.18%	0.646%
67	16.251	2217	2221	2227	rVB4	38971	55932	1.02%	0.065%
68	16.392	2240	2245	2252	rVB3	20462	49567	0.90%	0.057%
69	16.551	2267	2272	2277	rBV6	29109	46948	0.86%	0.054%
70	16.845	2318	2322	2326	rBV5	23370	37982	0.69%	0.044%
71	16.951	2336	2340	2344	rBV2	28940	37108	0.68%	0.043%
72	17.175	2375	2378	2380	rBV2	23601	25756	0.47%	0.030%
73	17.333	2402	2405	2413	rVB3	40209	63198	1.15%	0.073%
74	17.398	2413	2416	2420	rBV6	31913	45561	0.83%	0.053%
75	17.463	2420	2427	2439	rBV	2405633	3893755	70.99%	4.502%
76	17.563	2439	2444	2453	rBV2	2280315	4058073	73.99%	4.692%

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77	17.910	2501	2503	2509	rVB4	26385	31291	0.57%	0.036%
78	18.092	2531	2534	2539	rVB3	41831	53972	0.98%	0.062%
79	18.157	2542	2545	2549	rBV	80211	103211	1.88%	0.119%
80	18.204	2549	2553	2557	rVB5	29673	43190	0.79%	0.050%
81	18.257	2558	2562	2566	rBV	138428	198479	3.62%	0.229%
82	18.316	2567	2572	2581	rBV	1503743	2226716	40.60%	2.574%
83	18.957	2678	2681	2686	rVB3	57409	81067	1.48%	0.094%
84	19.186	2717	2720	2723	rBV	53282	52973	0.97%	0.061%
85	19.304	2737	2740	2746	rBV	79373	128357	2.34%	0.148%
86	19.369	2747	2751	2753	rBV	118095	138756	2.53%	0.160%
87	19.398	2753	2756	2758	rBV2	107569	121276	2.21%	0.140%
88	19.539	2776	2780	2788	rBV	450485	733897	13.38%	0.848%
89	19.792	2818	2823	2837	rVB	3241468	5484724	100.00%	6.341%
90	20.263	2900	2903	2907	rBV	163564	194010	3.54%	0.224%
91	20.745	2982	2985	2986	rBV	188633	197191	3.60%	0.228%
92	21.198	3059	3062	3064	rBV2	272083	320091	5.84%	0.370%
93	21.227	3064	3067	3073	rVB	749319	1041514	18.99%	1.204%
94	21.551	3118	3122	3135	rVB	2091323	3526586	64.30%	4.077%
95	22.121	3216	3219	3223	rVB	407282	458731	8.36%	0.530%
96	23.227	3401	3407	3414	rVB	1789170	2929240	53.41%	3.387%
97	23.927	3521	3526	3539	rVB2	2114072	4746970	86.55%	5.488%
98	24.098	3548	3555	3571	rVB	1708557	4271669	77.88%	4.938%
99	24.657	3643	3650	3661	rVB	2243071	4848456	88.40%	5.605%
100	26.609	3975	3982	3991	rBV	1019381	2668635	48.66%	3.085%

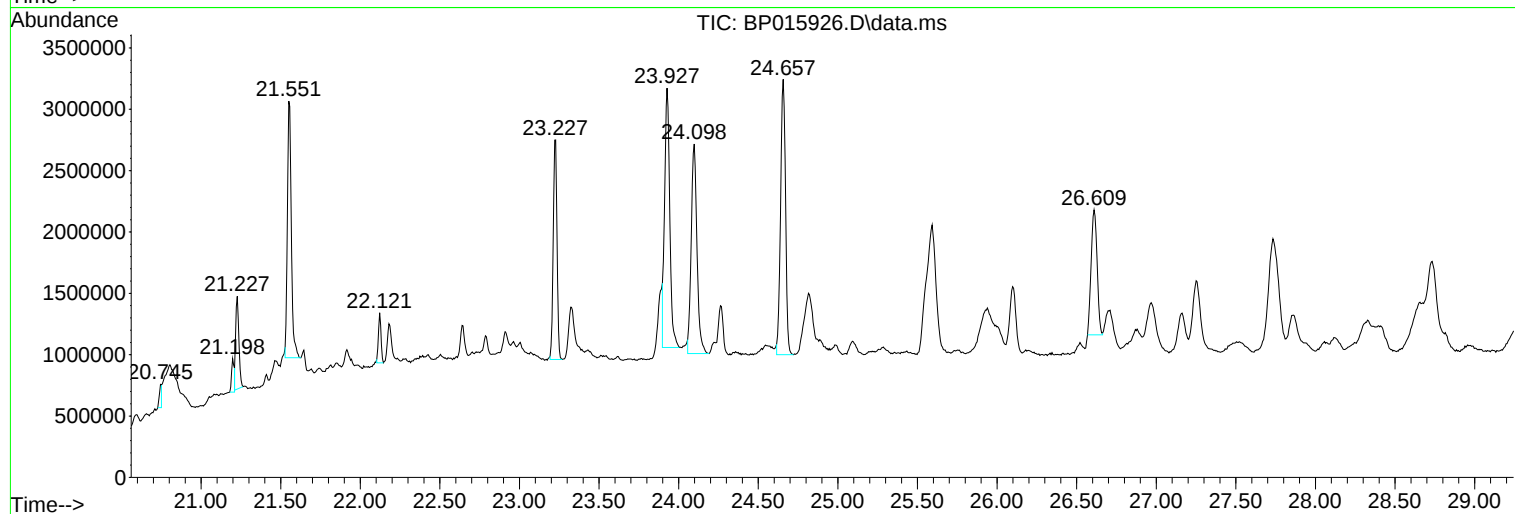
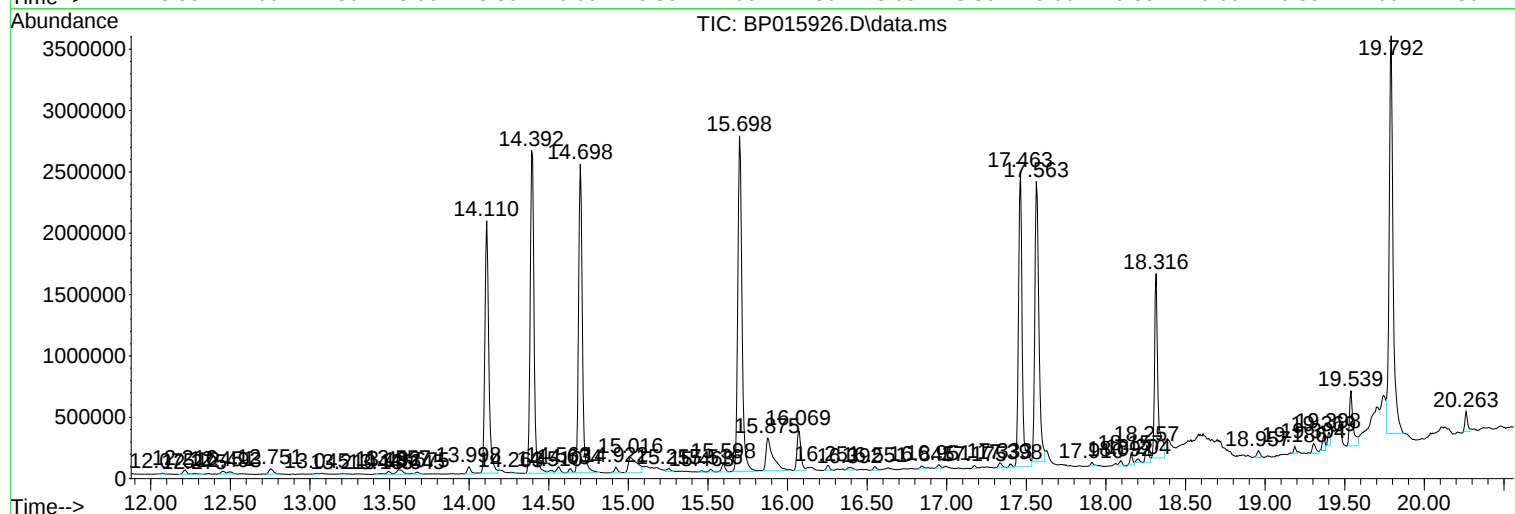
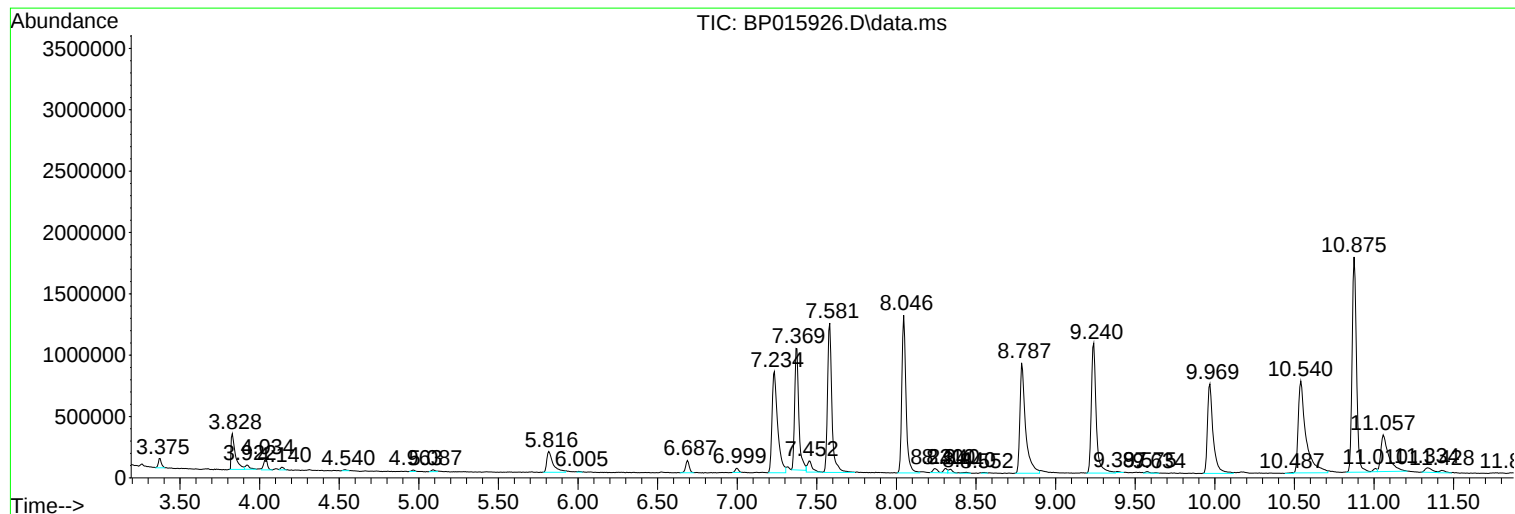
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
Quant Title : SVOA CALIBRATION

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



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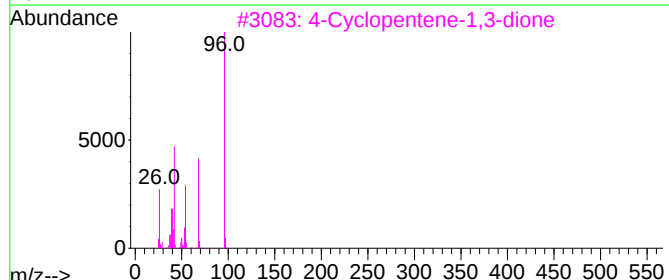
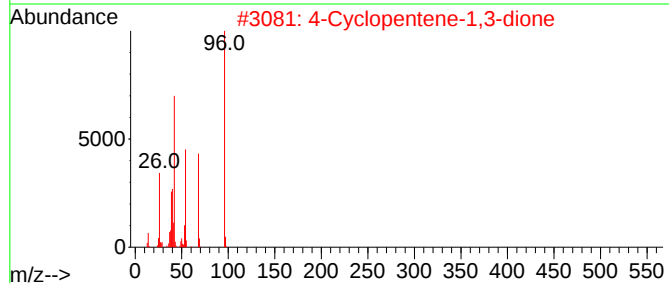
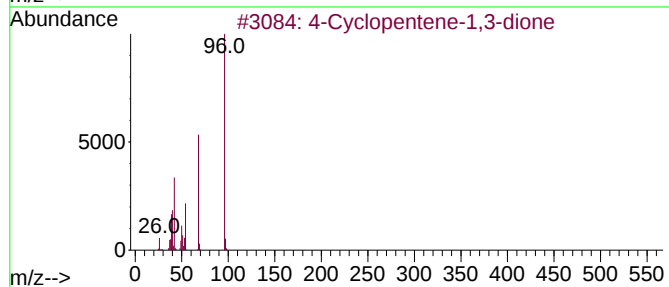
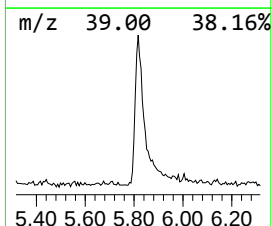
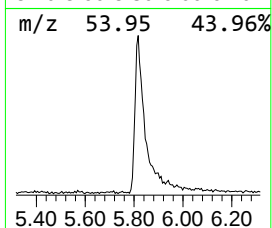
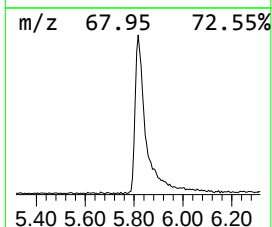
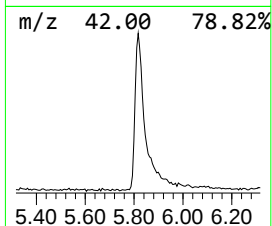
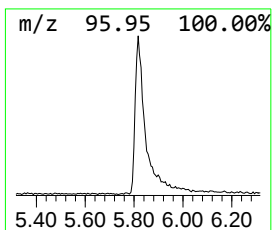
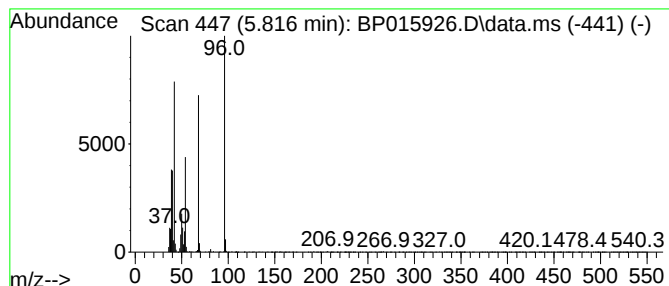
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 4-Cyclopentene-1,3-dione Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.816	4.06 ng/ul	475874	1,4-Dichlorobenzene-d4	8.046

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4-Cyclopentene-1,3-dione	96	C5H4O2	000930-60-9	95
2			4-Cyclopentene-1,3-dione	96	C5H4O2	000930-60-9	87
3			4-Cyclopentene-1,3-dione	96	C5H4O2	000930-60-9	83
4			3-Amino-1,2,4-triazine	96	C3H4N4	001120-99-6	47
5			But-1-ene-3-yne, 1-ethoxy-	96	C6H8O	002806-41-9	38



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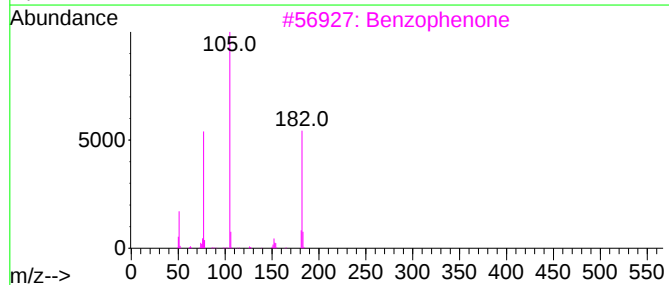
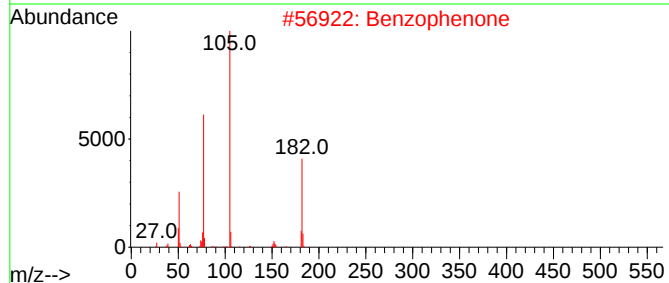
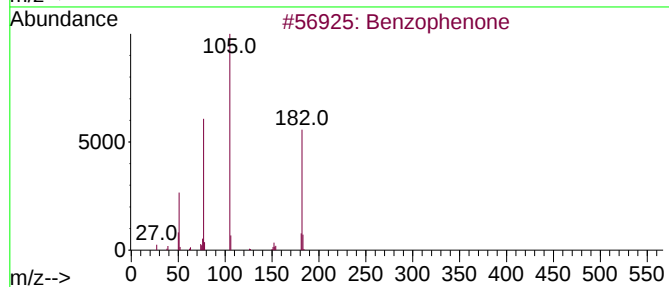
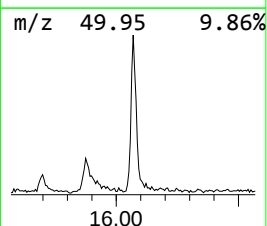
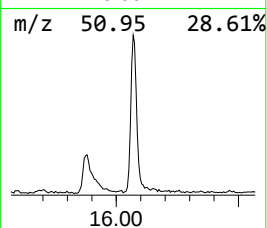
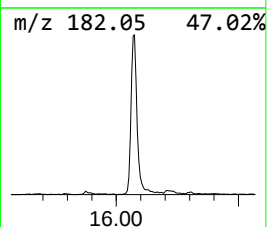
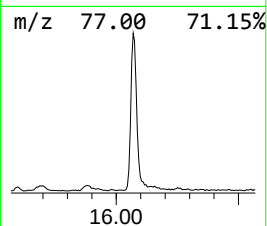
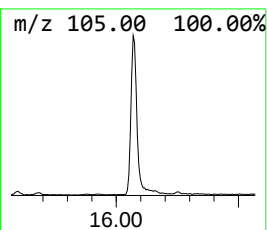
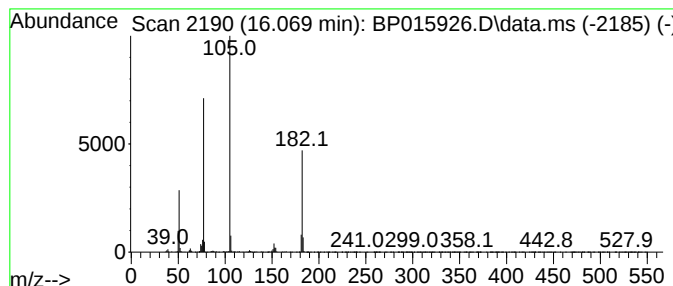
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Benzophenone Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.069	2.74 ng/ul	558475	Acenaphthene-d10	14.698

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



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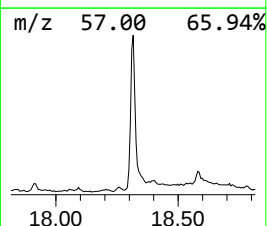
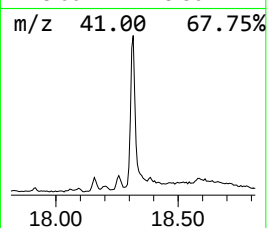
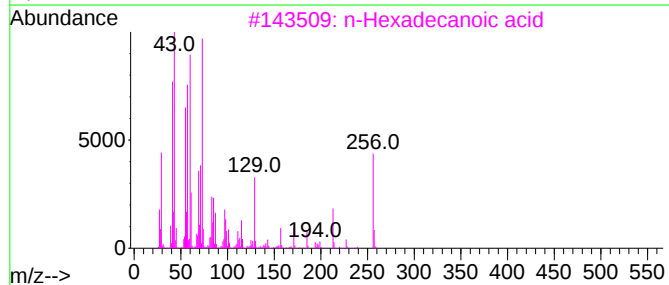
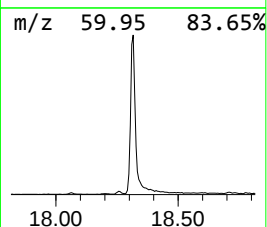
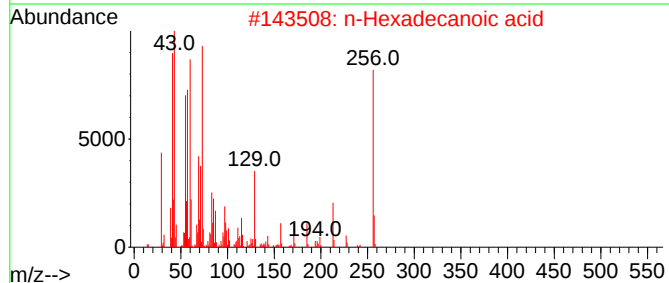
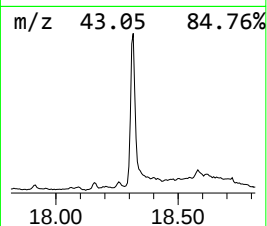
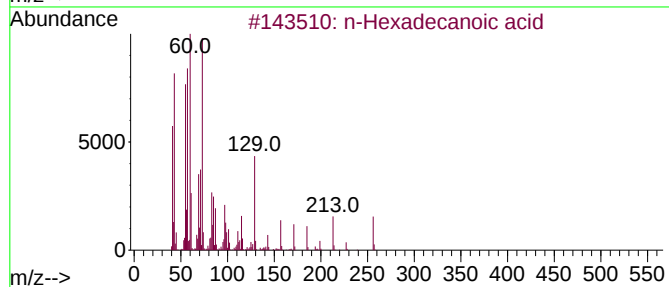
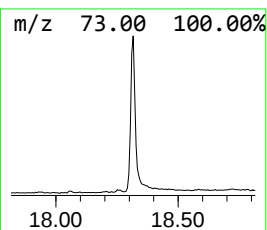
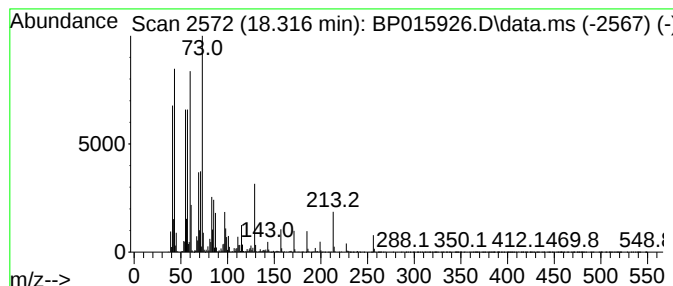
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 n-Hexadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.316	11.44 ng/ul	2226720	Phenanthrene-d10	17.463

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			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	97



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070223\
Data File : BP015926.D
Acq On : 02 Jul 2023 16:18
Operator : MA/JU
Sample : 03243-08
Misc :
ALS Vial : 91 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCGC8

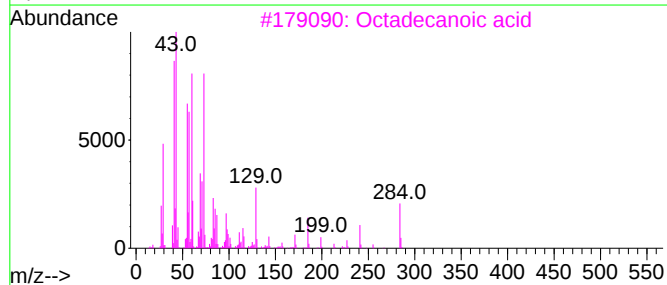
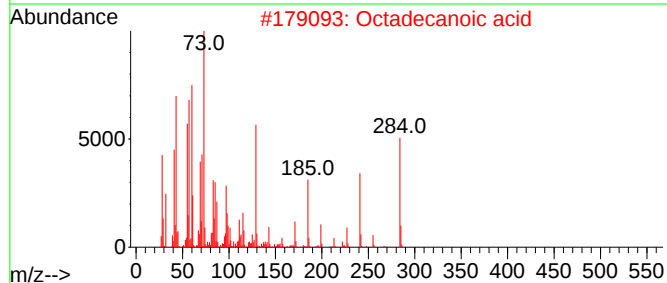
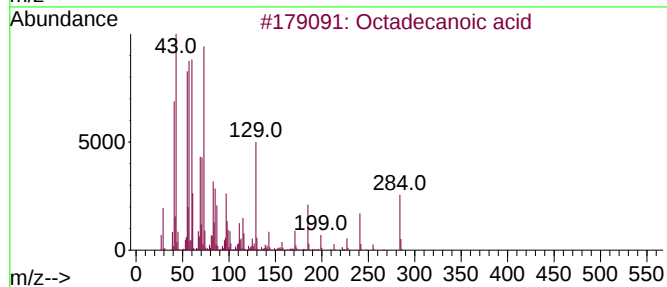
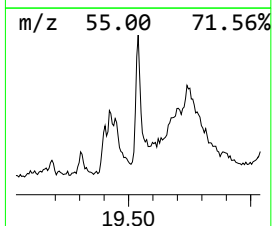
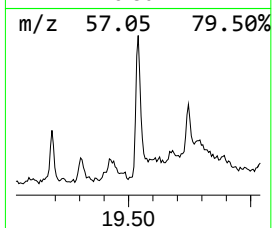
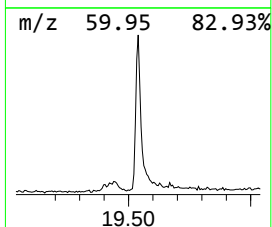
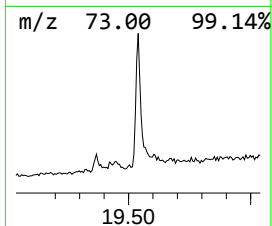
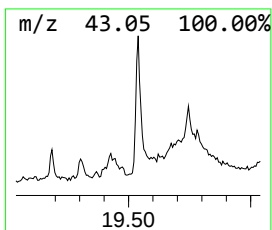
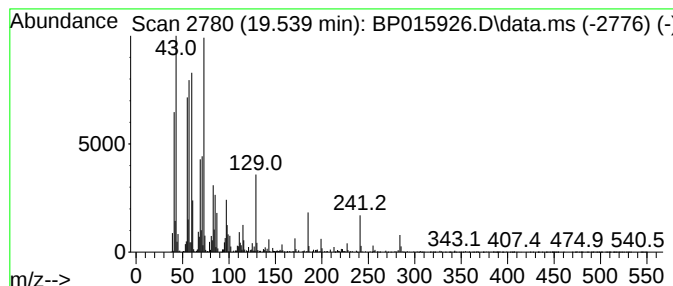
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Octadecanoic acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.539	4.16 ng/ul	733897	Chrysene-d12	21.551

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	96
4			Octadecanoic acid	284	C18H36O2	000057-11-4	95
5			Octadecanoic acid	284	C18H36O2	000057-11-4	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070223\
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ClientSampleId :
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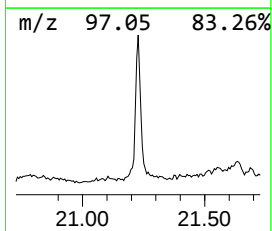
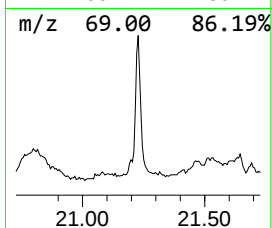
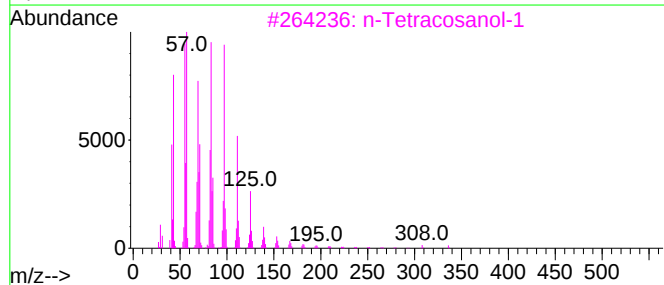
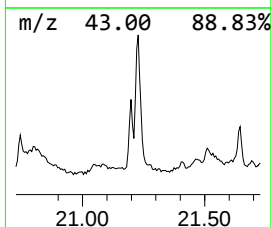
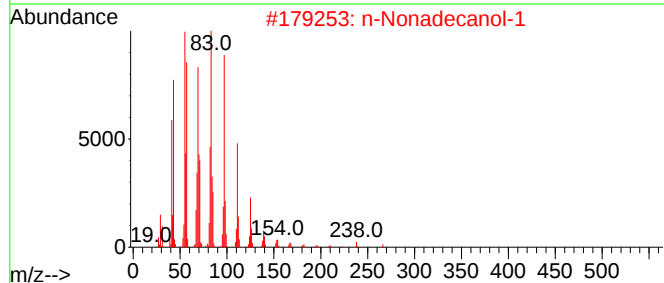
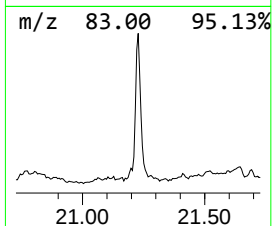
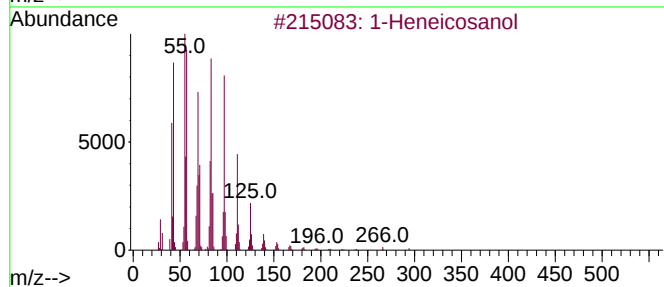
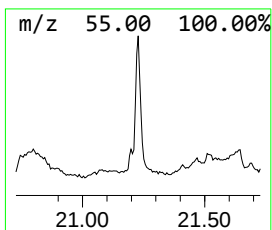
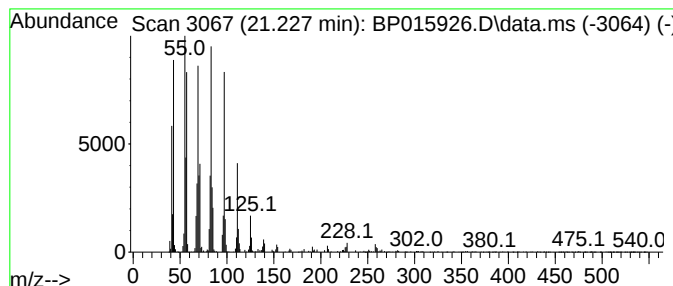
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 1-Heneicosanol Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.227	5.91 ng/ul	1041510	Chrysene-d12	21.551

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Heneicosanol	312	C21H44O	015594-90-8	93
2			n-Nonadecanol-1	284	C19H40O	001454-84-8	93
3			n-Tetracosanol-1	354	C24H50O	000506-51-4	93
4			1-Tetracosene	336	C24H48	010192-32-2	91
5			1-Octadecene	252	C18H36	000112-88-9	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070223\
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Operator : MA/JU
Sample : 03243-08
Misc :
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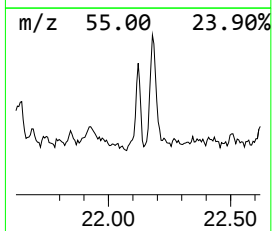
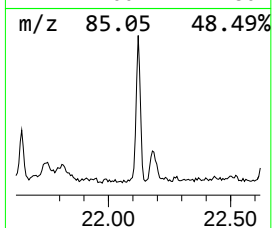
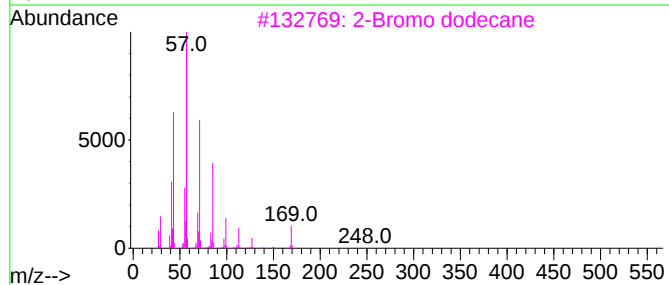
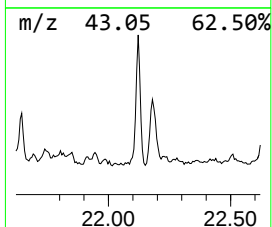
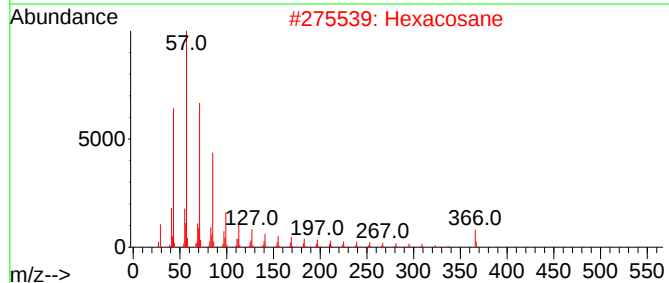
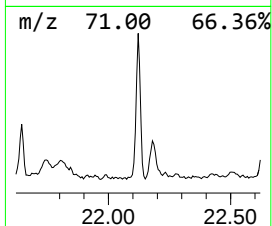
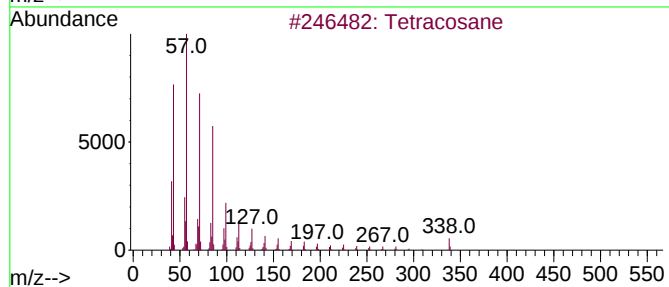
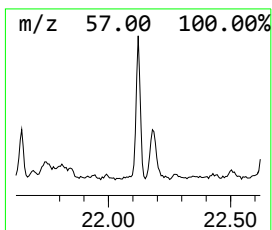
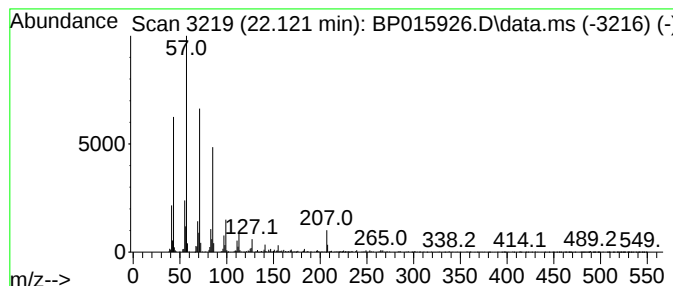
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD		R.T.	
22.121	2.60 ng/ul	458731	Chrysene-d12		21.551	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Tetracosane		338	C24H50	000646-31-1	95
2	Hexacosane		366	C26H54	000630-01-3	90
3	2-Bromo dodecane		248	C12H25Br	013187-99-0	90
4	Heneicosane		296	C21H44	000629-94-7	76
5	Docosane, 11-butyl-		366	C26H54	013475-76-8	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070223\
Data File : BP015926.D
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Sample : 03243-08
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ALS Vial : 91 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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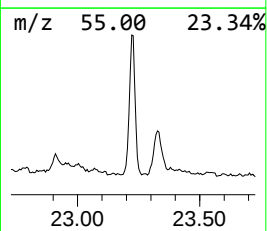
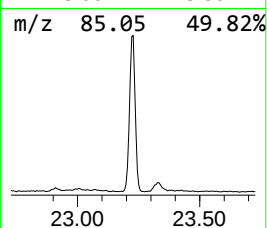
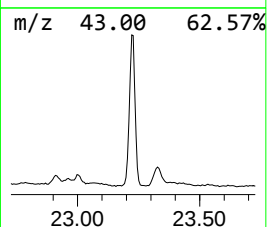
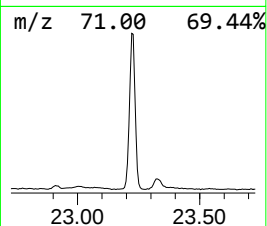
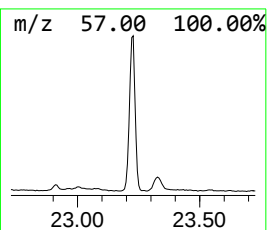
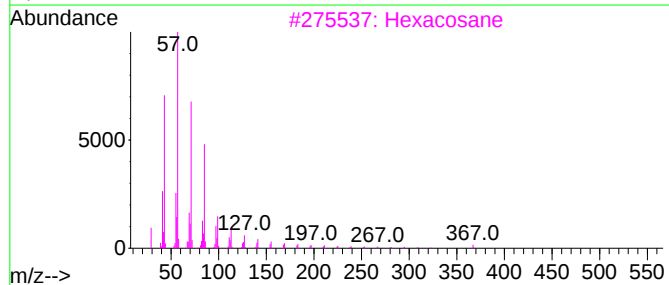
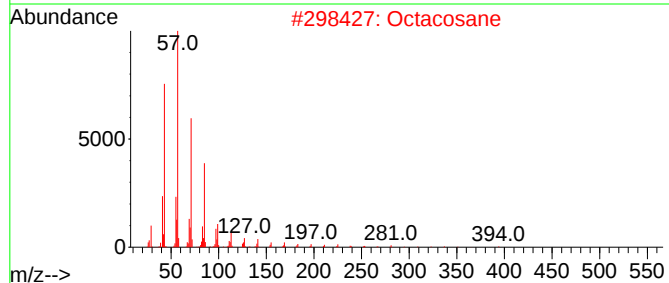
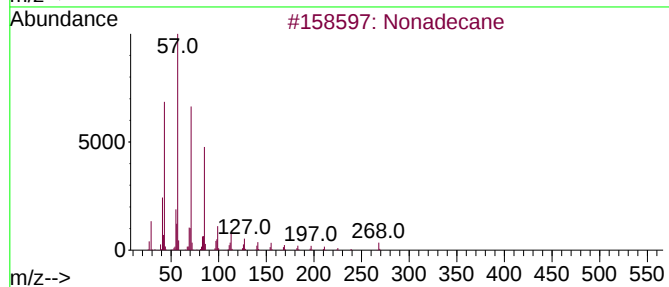
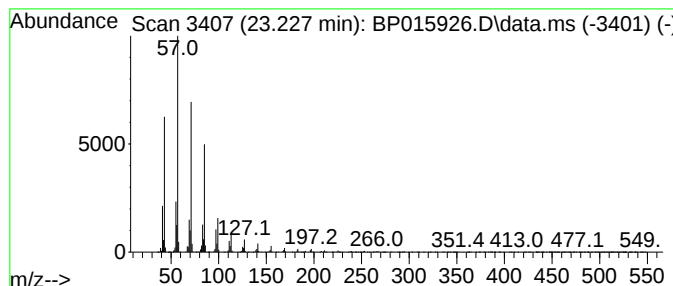
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.227	13.71 ng/ul	2929240	Perylene-d12	24.098

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Nonadecane	268	C19H40	000629-92-5	94
2		Octacosane	394	C28H58	000630-02-4	91
3		Hexacosane	366	C26H54	000630-01-3	91
4		Heneicosane	296	C21H44	000629-94-7	91
5		Dotriacontane, 1-iodo-	576	C32H65I	1000406-32-4	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070223\
Data File : BP015926.D
Acq On : 02 Jul 2023 16:18
Operator : MA/JU
Sample : 03243-08
Misc :
ALS Vial : 91 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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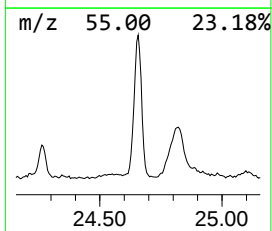
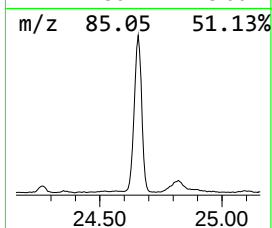
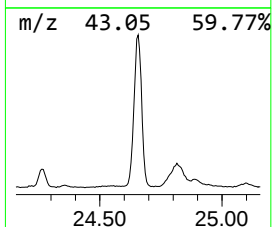
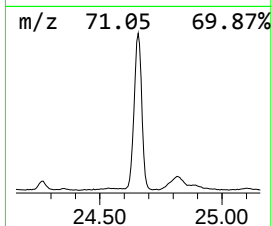
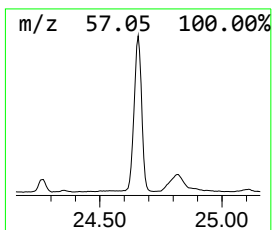
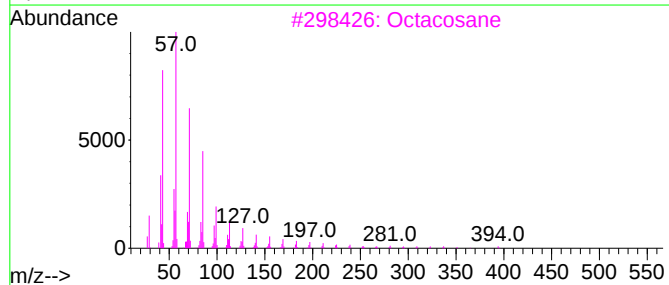
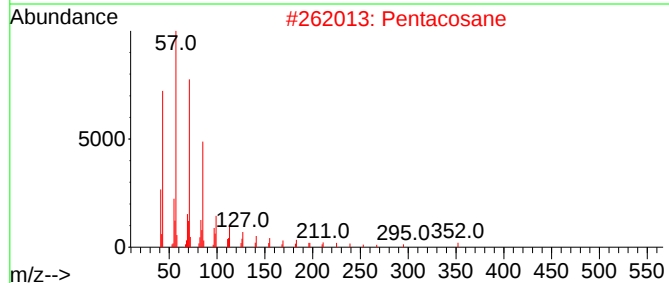
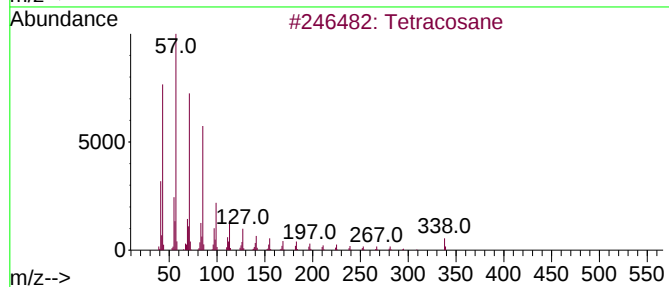
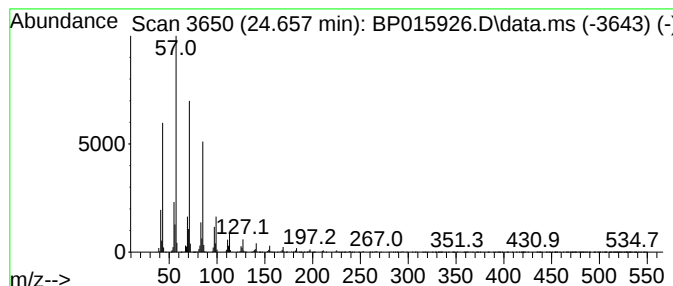
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.657	22.70 ng/ul	4848460	Perylene-d12	24.098

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetracosane	338	C24H50	000646-31-1	96
2			Pentacosane	352	C25H52	000629-99-2	94
3			Octacosane	394	C28H58	000630-02-4	91
4			Heneicosane	296	C21H44	000629-94-7	91
5			Heneicosane	296	C21H44	000629-94-7	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070223\
Data File : BP015926.D
Acq On : 02 Jul 2023 16:18
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Sample : 03243-08
Misc :
ALS Vial : 91 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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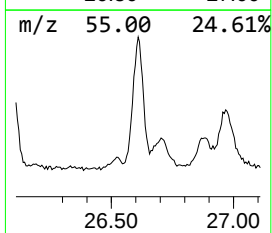
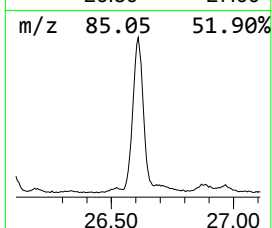
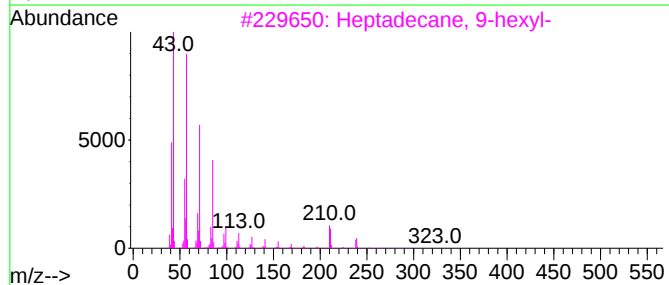
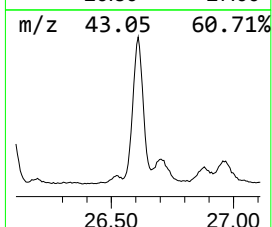
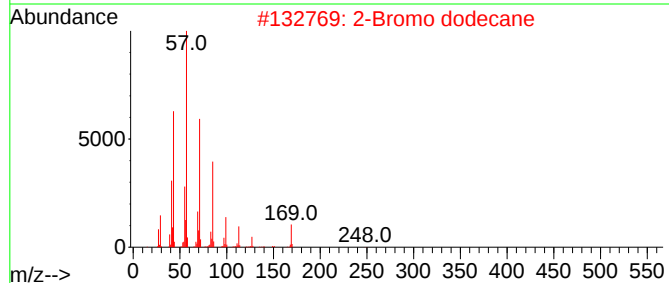
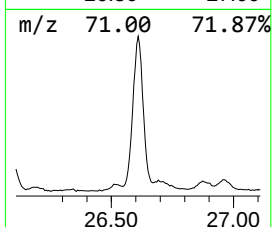
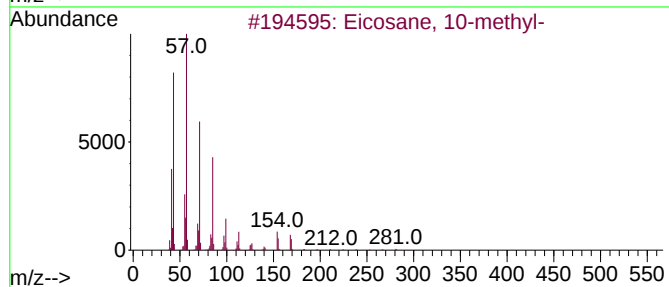
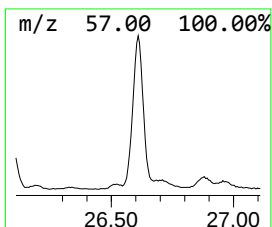
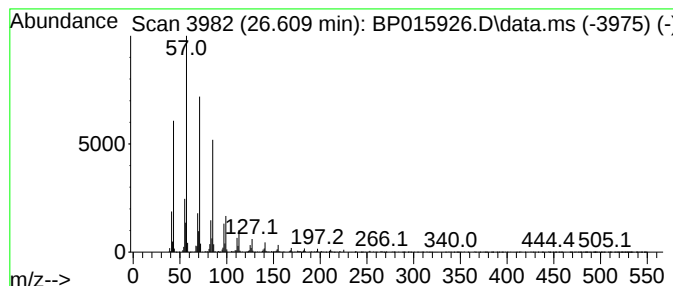
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.609	12.49 ng/ul	2668640	Perylene-d12	24.098

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Eicosane, 10-methyl-	296	C21H44	054833-23-7	94
2		2-Bromo dodecane	248	C12H25Br	013187-99-0	93
3		Heptadecane, 9-hexyl-	324	C23H48	055124-79-3	93
4		Heneicosane, 11-(1-ethylpropyl)-	366	C26H54	055282-11-6	91
5		Hexacosane	366	C26H54	000630-01-3	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070223\
Data File : BP015926.D
Acq On : 02 Jul 2023 16:18
Operator : MA/JU
Sample : 03243-08
Misc :
ALS Vial : 91 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCGC8

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.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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Benzophenone	16.069	2.7	ng/ul	558475	3	14.698	4069300	20.0
n-Hexadecanoic ...	18.316	11.4	ng/ul	2226720	4	17.463	3893760	20.0
Octadecanoic acid	19.539	4.2	ng/ul	733897	5	21.551	3526590	20.0
1-Heneicosanol	21.227	5.9	ng/ul	1041510	5	21.551	3526590	20.0
(DEL) Alkane: S...	22.121	2.6	ng/ul	458731	5	21.551	3526590	20.0
(DEL) Alkane: S...	23.227	13.7	ng/ul	2929240	6	24.098	4271670	20.0
(DEL) Alkane: S...	24.657	22.7	ng/ul	4848460	6	24.098	4271670	20.0
(DEL) Alkane: S...	26.609	12.5	ng/ul	2668640	6	24.098	4271670	20.0