

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM051520\
 Data File : BM025988.D
 Acq On : 15 May 2020 14:45
 Operator : MA/SJ
 Sample : L2660-04
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 CB7C3

Integration Parameters: LSCINT.P

Integrator: RTE
 Smoothing : OFF
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0

Filtering: 5
 Min Area: 0.1 % of largest Peak
 Max Peaks: 100
 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
 Title : SVOA CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.081	29	33	43	rVB	46275	74131	1.80%	0.159%
2	3.787	149	153	157	rBV3	7821	10289	0.25%	0.022%
3	4.687	301	306	313	rVB7	3366	7382	0.18%	0.016%
4	5.640	464	468	481	rBV	85947	145657	3.53%	0.312%
5	6.140	549	553	557	rBV4	7066	14719	0.36%	0.031%
6	6.193	559	562	568	rVB6	6038	10088	0.24%	0.022%
7	6.693	641	647	662	rVB	785824	1246669	30.21%	2.667%
8	6.857	667	675	688	rBV	614948	946573	22.94%	2.025%
9	7.045	699	707	719	rBV	782789	1177430	28.54%	2.519%
10	7.504	778	785	793	rBV	642505	992651	24.06%	2.124%
11	7.593	795	800	807	rVB2	15318	23596	0.57%	0.050%
12	8.216	898	906	917	rBV	952631	1425773	34.56%	3.050%
13	8.651	972	980	991	rBV	775736	1214243	29.43%	2.598%
14	8.851	1008	1014	1022	rVB3	12433	25305	0.61%	0.054%
15	9.128	1056	1061	1065	rVB	26499	39152	0.95%	0.084%
16	9.363	1094	1101	1111	rBV	545843	862943	20.91%	1.846%
17	9.898	1184	1192	1204	rBV	855451	1367473	33.14%	2.925%
18	10.269	1246	1255	1264	rBV	835274	1330803	32.25%	2.847%
19	10.410	1272	1279	1297	rBV	653298	1107204	26.83%	2.369%
20	11.875	1523	1528	1534	rBV	12789	20816	0.50%	0.045%
21	12.510	1630	1636	1640	rBV9	6042	11656	0.28%	0.025%
22	12.710	1666	1670	1674	rVB6	2375	4754	0.12%	0.010%
23	12.763	1674	1679	1684	rBV5	6611	11063	0.27%	0.024%
24	13.004	1716	1720	1724	rBV3	9575	13193	0.32%	0.028%
25	13.069	1724	1731	1737	rBV5	7227	11291	0.27%	0.024%
26	13.574	1810	1817	1821	rBV	1404807	1986800	48.15%	4.250%
27	13.616	1821	1824	1833	rVB	454388	595044	14.42%	1.273%
28	13.839	1855	1862	1873	rBV	1730353	2460963	59.64%	5.265%
29	14.086	1899	1904	1908	rBV2	10456	15937	0.39%	0.034%
30	14.151	1908	1915	1923	rVV	1147916	1604092	38.88%	3.432%
31	14.363	1945	1951	1967	rBV	756331	1045096	25.33%	2.236%
32	14.986	2052	2057	2061	rBV6	4089	6889	0.17%	0.015%
33	15.045	2061	2067	2071	rVV4	6809	9458	0.23%	0.020%
34	15.151	2078	2085	2093	rBV	2306260	3082916	74.72%	6.595%

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Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
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35	15.274	2100	2106	2119	rBV	617513	851899	20.65%	1.822%
36	15.392	2122	2126	2130	rVB2	25123	34050	0.83%	0.073%
37	15.604	2159	2162	2168	rBV7	3448	6544	0.16%	0.014%
38	15.910	2211	2214	2216	rBV3	4710	6044	0.15%	0.013%
39	15.939	2216	2219	2222	rVV4	9254	13979	0.34%	0.030%
40	16.163	2254	2257	2263	rVB6	2936	4675	0.11%	0.010%
41	16.586	2327	2329	2333	rVB3	3610	4269	0.10%	0.009%
42	16.692	2342	2347	2352	rVB4	5852	11415	0.28%	0.024%
43	16.898	2376	2382	2388	rBV	1367416	1834832	44.47%	3.925%
44	16.998	2392	2399	2414	rVB	2462137	3313013	80.30%	7.087%
45	17.315	2449	2453	2458	rVB9	2878	5509	0.13%	0.012%
46	17.827	2535	2540	2545	rBV	110933	158370	3.84%	0.339%
47	18.127	2586	2591	2595	rBV6	6435	8498	0.21%	0.018%
48	18.327	2622	2625	2629	rBV6	2560	4460	0.11%	0.010%
49	18.633	2675	2677	2684	rBV8	4097	6647	0.16%	0.014%
50	18.768	2697	2700	2706	rVB6	8686	10773	0.26%	0.023%
51	18.933	2723	2728	2731	rBV	25067	30930	0.75%	0.066%
52	19.115	2755	2759	2766	rBV	50774	81302	1.97%	0.174%
53	19.186	2767	2771	2778	rVB	24617	39277	0.95%	0.084%
54	19.298	2784	2790	2797	rBV	3053386	3907682	94.71%	8.360%
55	19.356	2797	2800	2803	rVV3	32127	37973	0.92%	0.081%
56	19.392	2803	2806	2811	rVB6	13183	17792	0.43%	0.038%
57	19.856	2882	2885	2888	rBV5	12614	17192	0.42%	0.037%
58	19.892	2888	2891	2897	rVV2	29867	35242	0.85%	0.075%
59	20.392	2972	2976	2980	rBV	57308	57864	1.40%	0.124%
60	20.845	3049	3053	3060	rBV2	522926	708356	17.17%	1.515%
61	21.098	3090	3096	3101	rBV	1765808	2133024	51.70%	4.563%
62	21.292	3125	3129	3137	rBV	277928	299485	7.26%	0.641%
63	21.709	3196	3200	3207	rBV	432911	521398	12.64%	1.115%
64	22.150	3271	3275	3280	rBV	552842	656897	15.92%	1.405%
65	22.633	3352	3357	3362	rBV	553843	737547	17.88%	1.578%
66	23.092	3429	3435	3442	rBV	2488164	4126045	100.00%	8.827%
67	23.162	3442	3447	3451	rVV	528338	788488	19.11%	1.687%
68	23.221	3451	3457	3470	rVB2	1288054	2306485	55.90%	4.934%
69	23.762	3544	3549	3556	rBV	371879	633750	15.36%	1.356%
70	24.456	3662	3667	3681	rVB	219799	430856	10.44%	0.922%

LSC Area Percent Report

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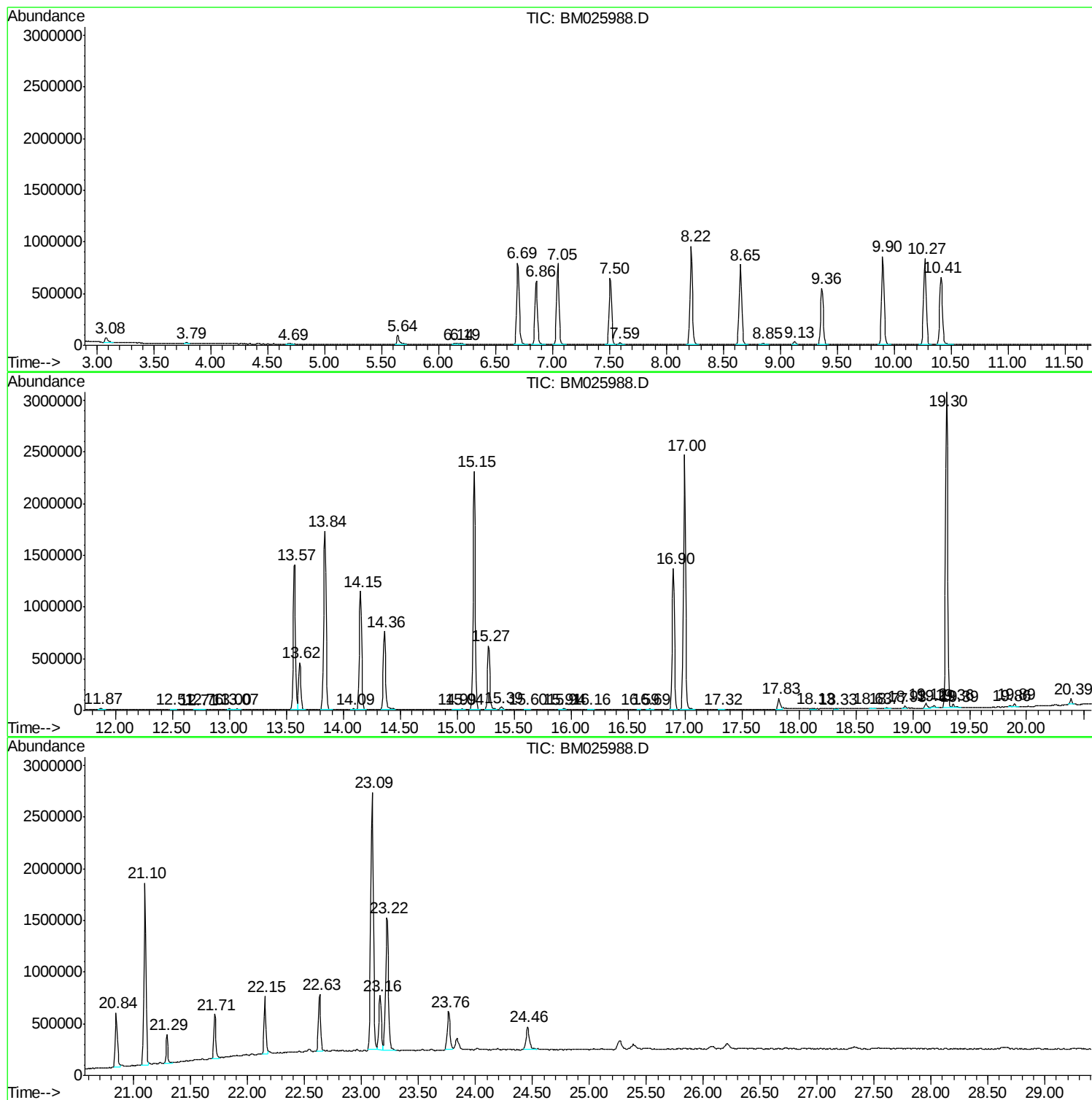
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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Sample : L2660-04
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Instrument :
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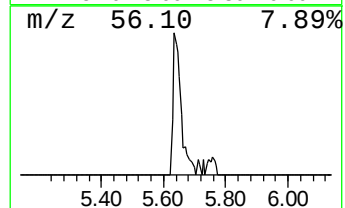
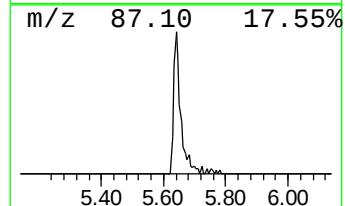
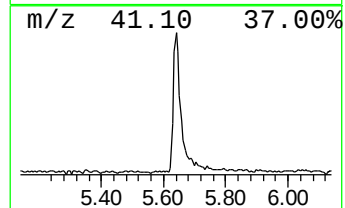
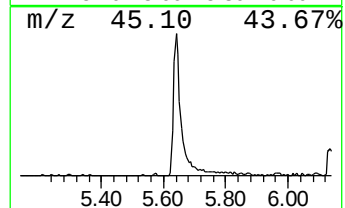
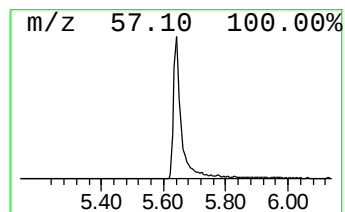
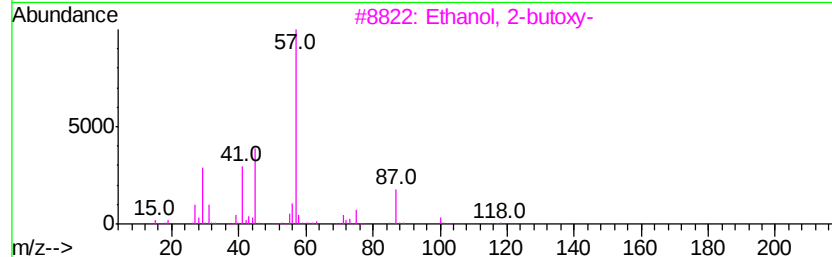
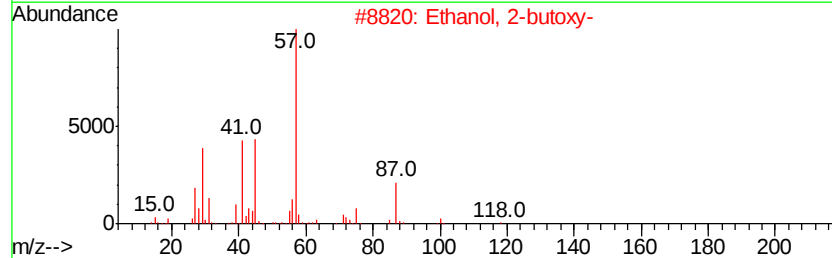
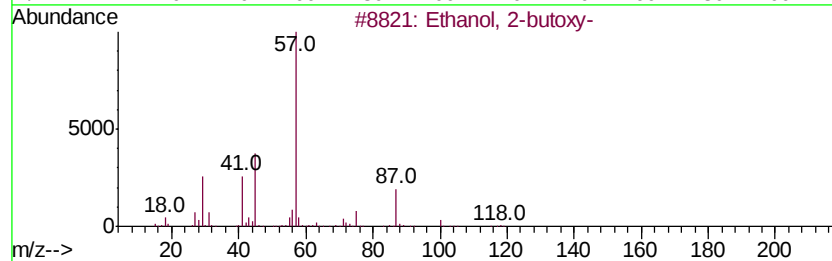
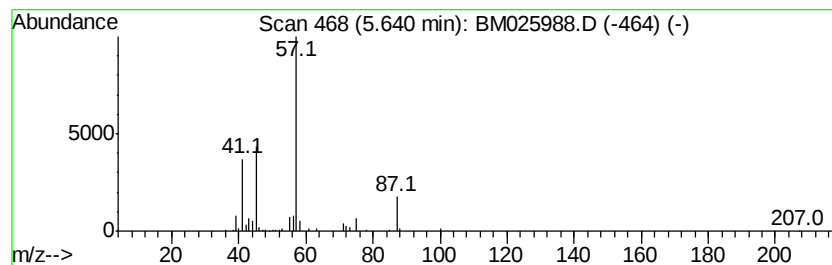
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Ethanol, 2-butoxy- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.64	2.93 ng/ul	145657	1,4-Dichlorobenzene-d4	7.50

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Ethanol, 2-butoxy-	118	C6H14O2	000111-76-2	78
2		Ethanol, 2-butoxy-	118	C6H14O2	000111-76-2	64
3		Ethanol, 2-butoxy-	118	C6H14O2	000111-76-2	53
4		Ethylene glycol monoisobutyl ether	118	C6H14O2	004439-24-1	45
5		3,3-Dimethylbutane-2-ol	102	C6H14O	000464-07-3	36



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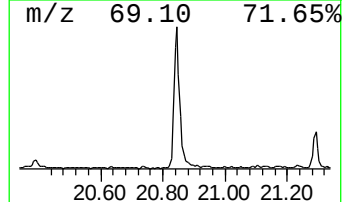
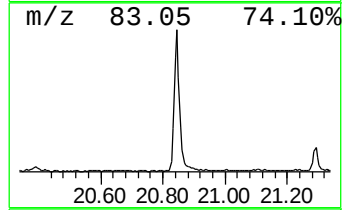
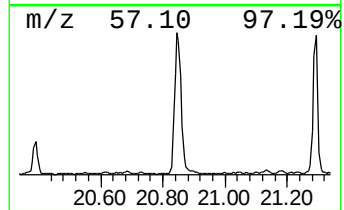
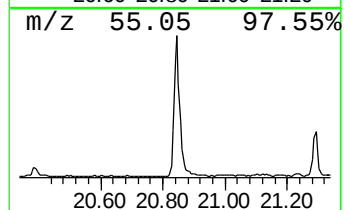
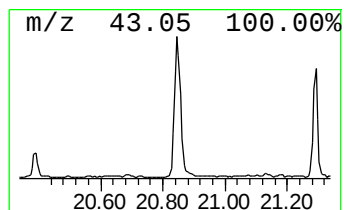
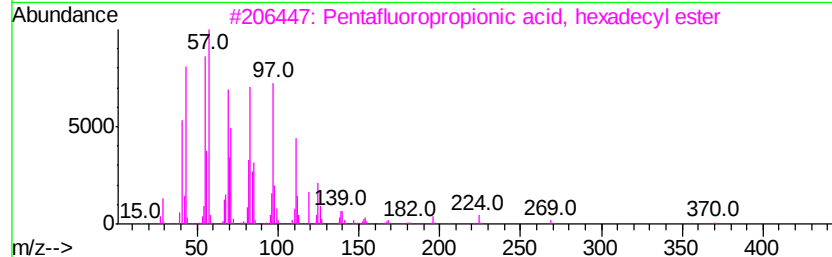
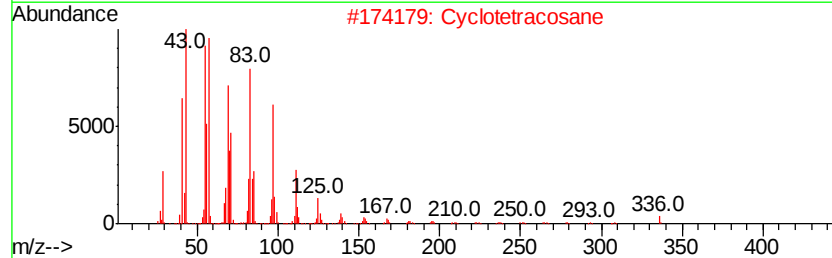
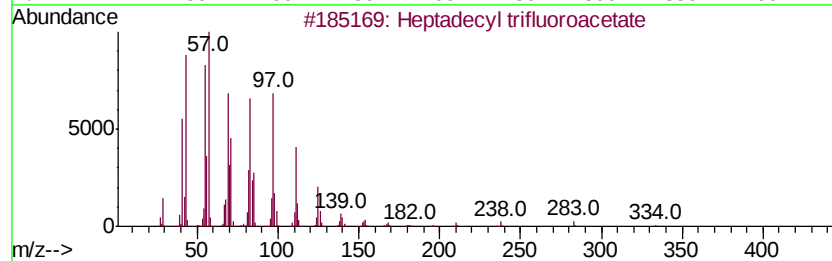
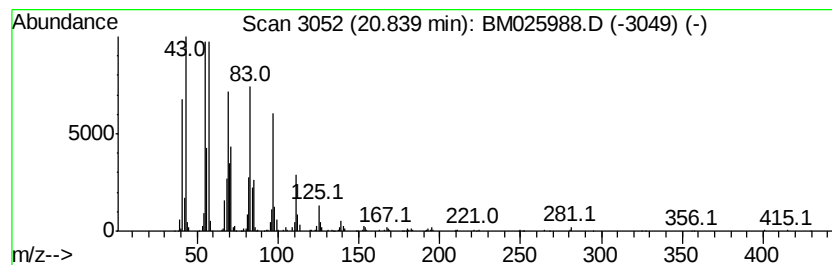
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.84	6.64 ng/ul	708356	Chrysene-d12	21.10

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptadecyl trifluoroacetate	352	C19H35F3O2	1000351-87-0	91
2		Cyclotetracosane	336	C24H48	000297-03-0	91
3		Pentafluoropropionic acid, hexadecyl ester	388	C19H33F5O2	006222-07-7	91
4		Cyclooctacosane	392	C28H56	000297-24-5	91
5		1-Tricosene	322	C23H46	018835-32-0	91



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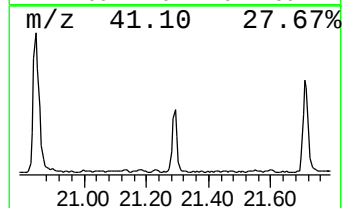
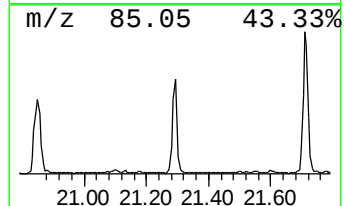
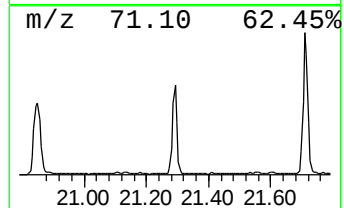
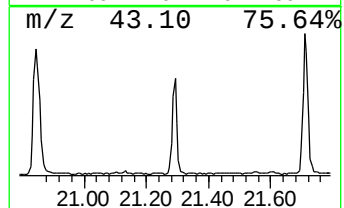
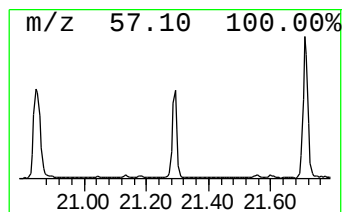
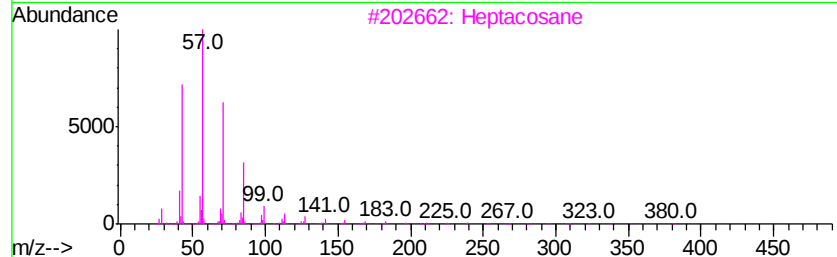
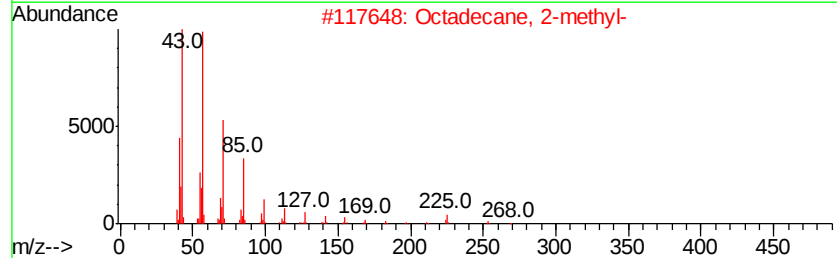
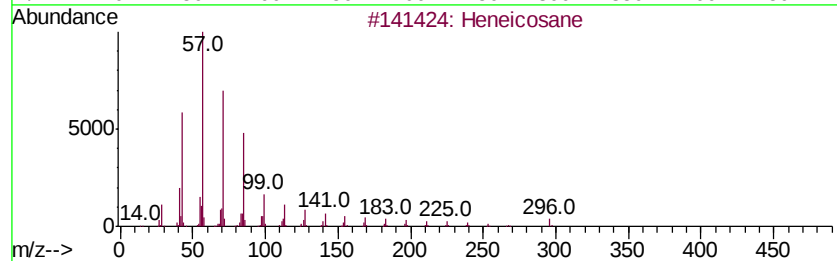
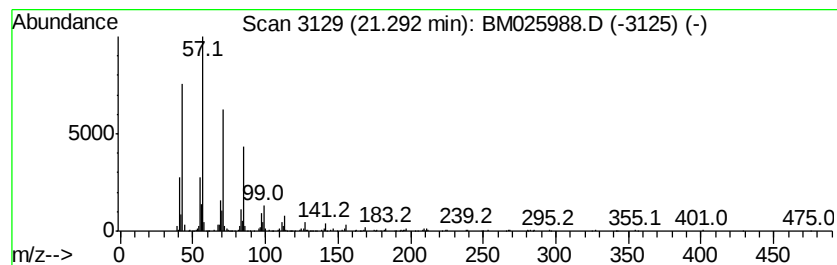
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.29	2.81 ng/ul	299485	Chrysene-d12	21.10

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosane	296	C21H44	000629-94-7	95
2		Octadecane, 2-methyl-	268	C19H40	001560-88-9	93
3		Heptacosane	380	C27H56	000593-49-7	91
4		Eicosane, 7-hexyl-	366	C26H54	055333-99-8	91
5		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91



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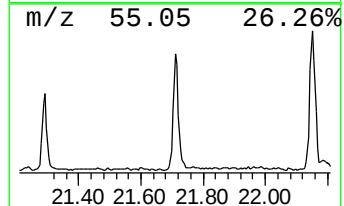
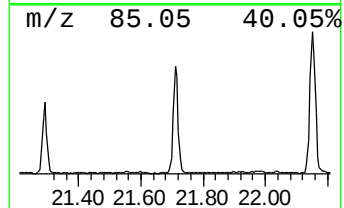
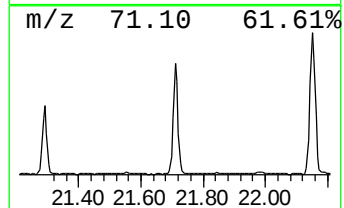
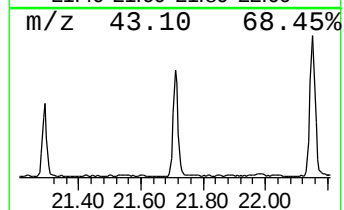
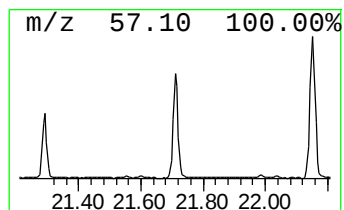
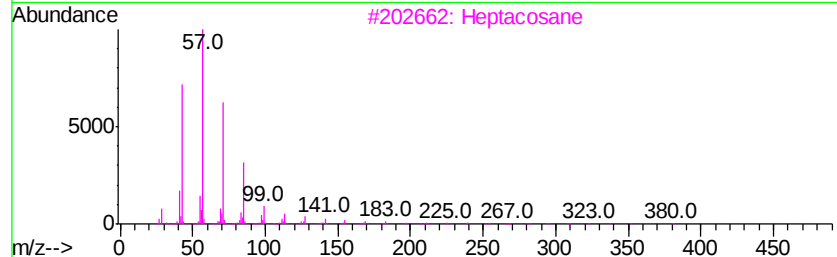
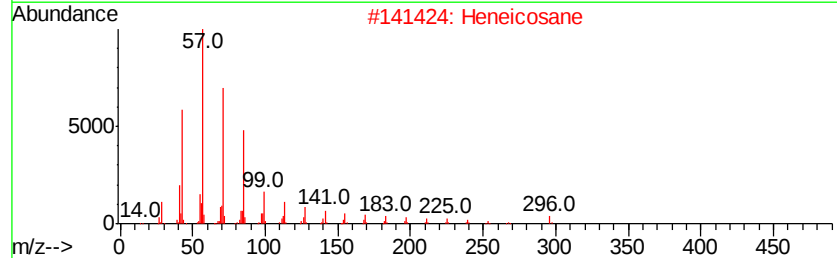
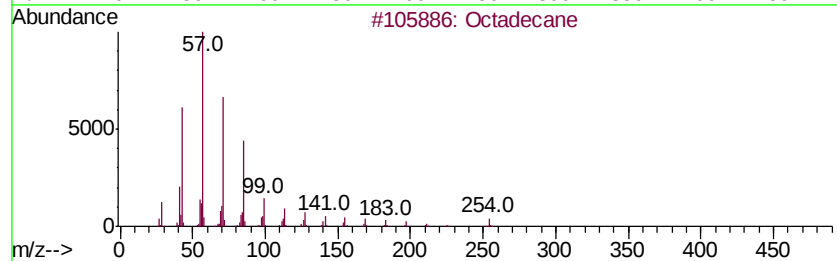
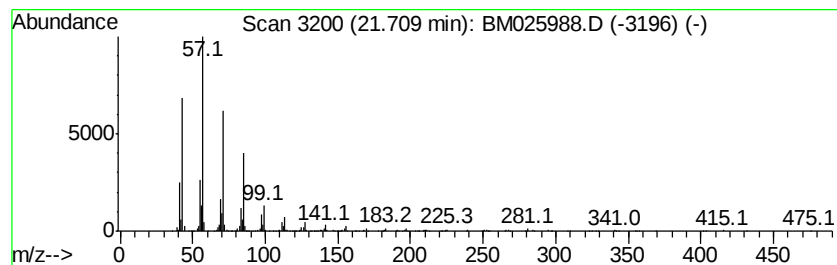
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.71	4.89 ng/ul	521398	Chrysene-d12	21.10

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecane	254	C18H38	000593-45-3	95
2			Heneicosane	296	C21H44	000629-94-7	95
3			Heptacosane	380	C27H56	000593-49-7	91
4			Hentriacontane	437	C31H64	000630-04-6	91
5			Tetratetracontane	619	C44H90	007098-22-8	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM051520\
Data File : BM025988.D
Acq On : 15 May 2020 14:45
Operator : MA/SJ
Sample : L2660-04
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
CB7C3

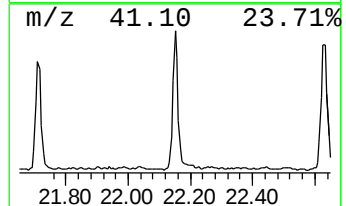
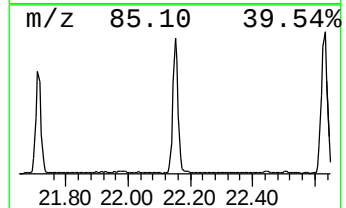
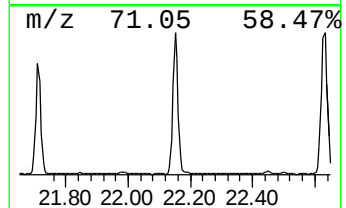
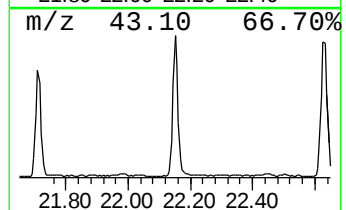
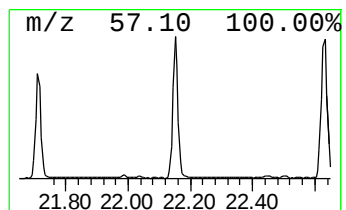
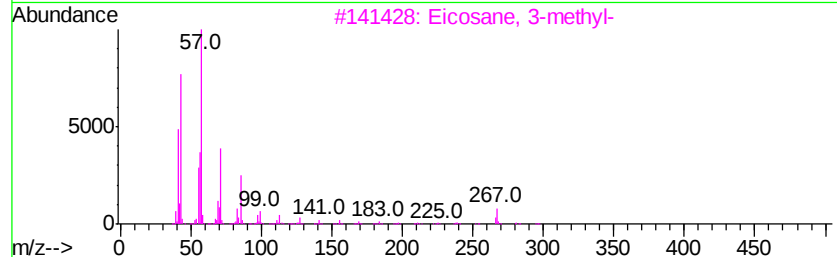
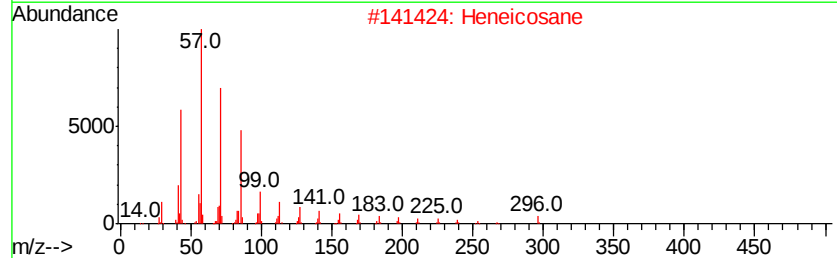
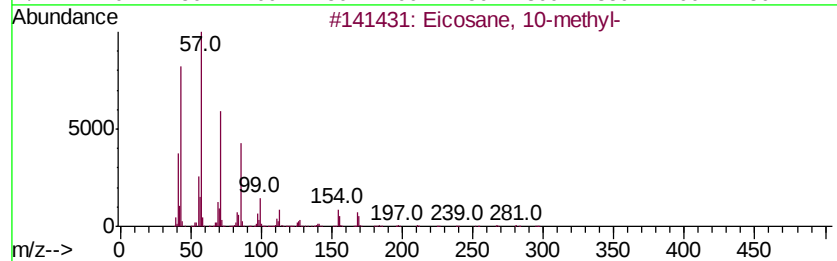
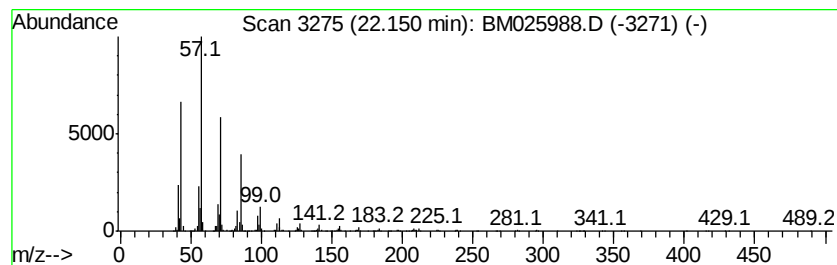
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.15	6.16 ng/ul	656897	Chrysene-d12	21.10

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane, 10-methyl-		296	C21H44	054833-23-7	97
2	Heneicosane		296	C21H44	000629-94-7	95
3	Eicosane, 3-methyl-		296	C21H44	006418-46-8	95
4	Heneicosane		296	C21H44	000629-94-7	95
5	1-Iodo-2-methylundecane		296	C12H25I	073105-67-6	94



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM051520\
Data File : BM025988.D
Acq On : 15 May 2020 14:45
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Sample : L2660-04
Misc :
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Instrument :
BNA_M
ClientSampleId :
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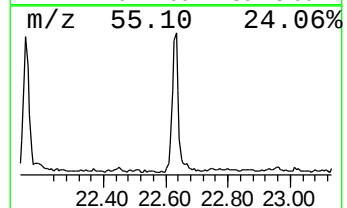
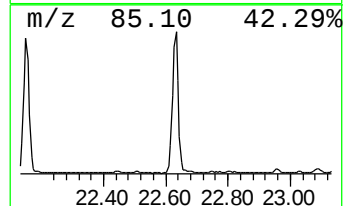
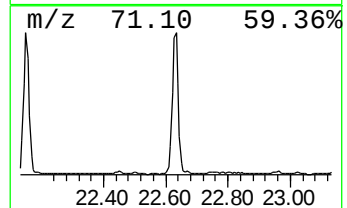
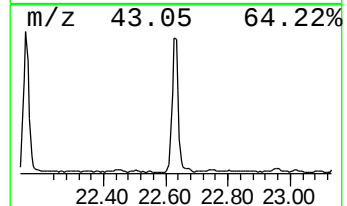
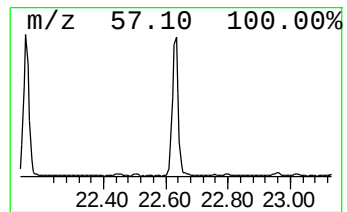
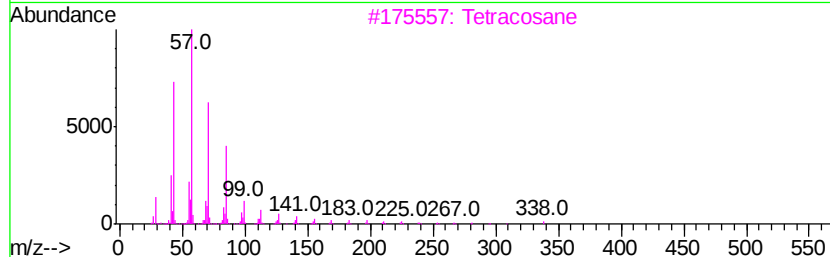
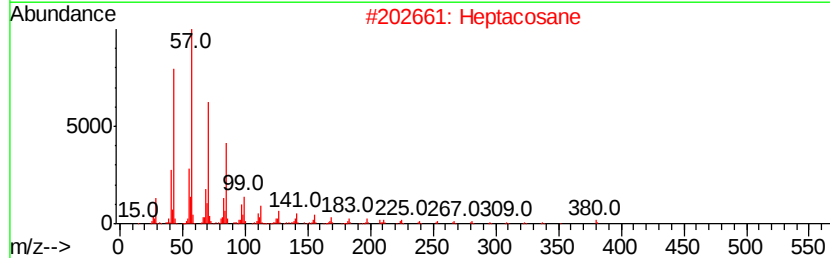
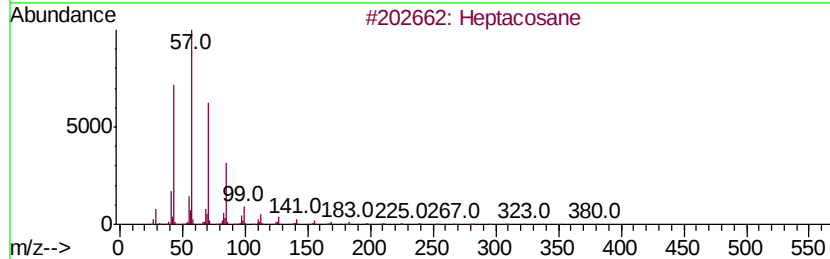
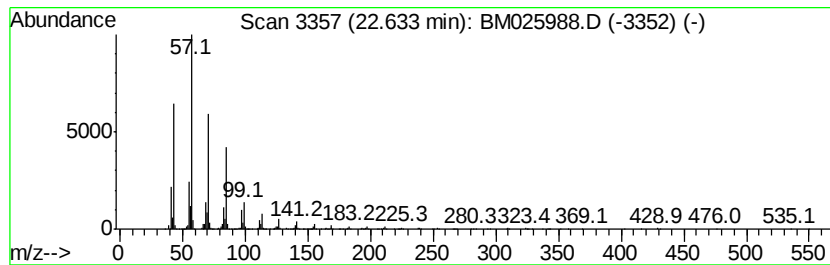
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.63	6.40 ng/ul	737547	Perylene-d12	23.22

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptacosane	380	C27H56	000593-49-7	91
2		Heptacosane	380	C27H56	000593-49-7	91
3		Tetracosane	338	C24H50	000646-31-1	91
4		Hexatriacontane	507	C36H74	000630-06-8	91
5		Heneicosane	296	C21H44	000629-94-7	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM051520\
Data File : BM025988.D
Acq On : 15 May 2020 14:45
Operator : MA/SJ
Sample : L2660-04
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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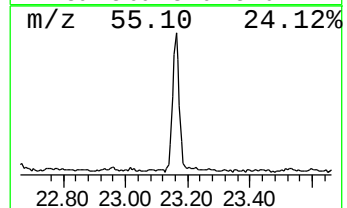
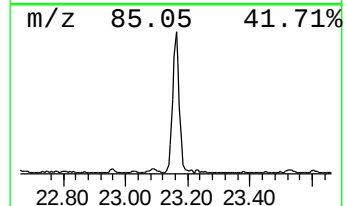
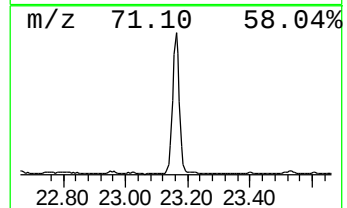
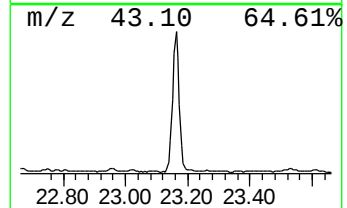
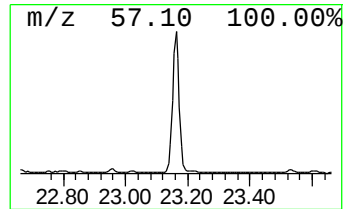
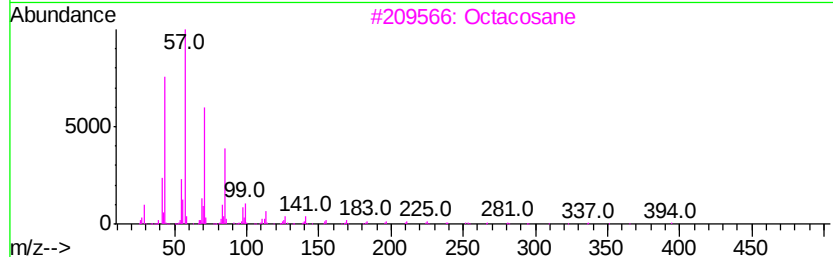
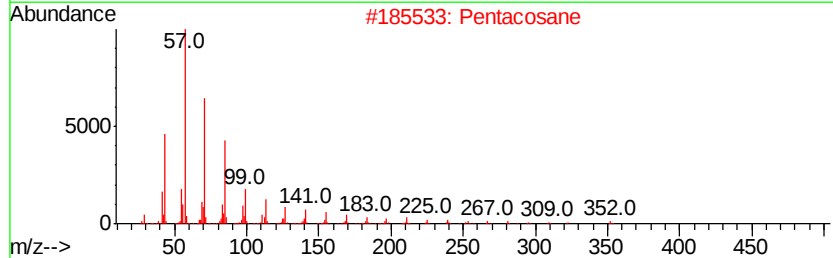
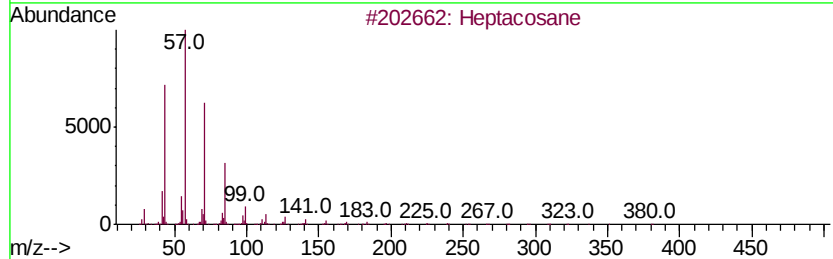
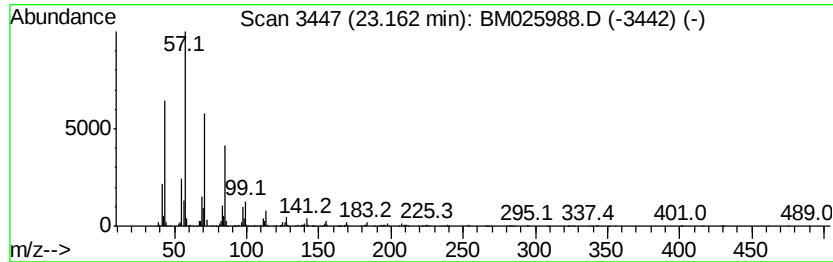
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.16	6.84 ng/ul	788488	Perylene-d12	23.22

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptacosane	380	C27H56	000593-49-7	91
2			Pentacosane	352	C25H52	000629-99-2	91
3			Octacosane	394	C28H58	000630-02-4	91
4			Hexacosane	366	C26H54	000630-01-3	91
5			Octacosane	394	C28H58	000630-02-4	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM051520\
Data File : BM025988.D
Acq On : 15 May 2020 14:45
Operator : MA/SJ
Sample : L2660-04
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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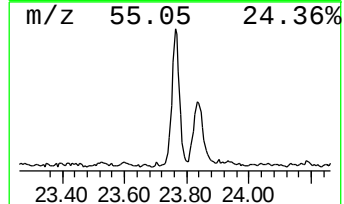
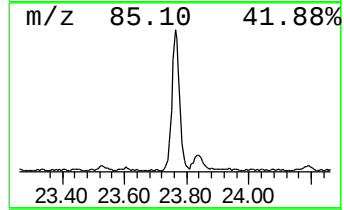
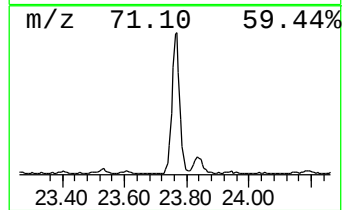
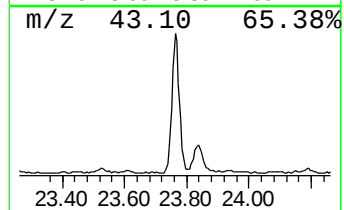
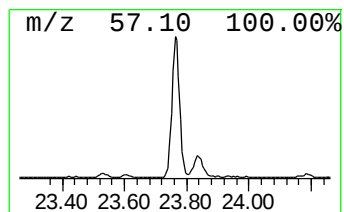
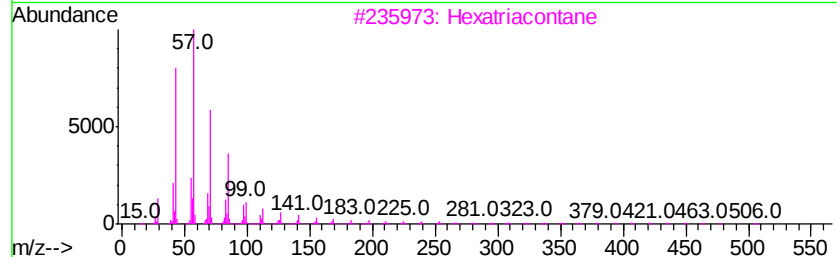
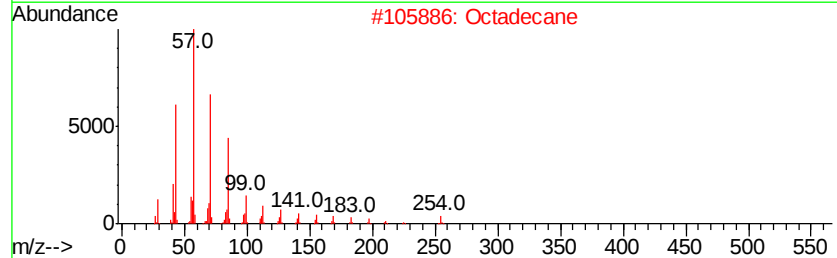
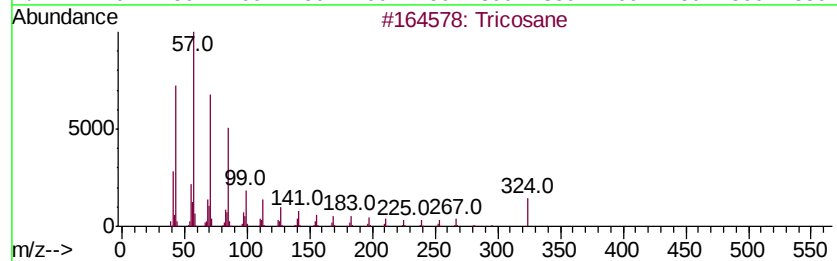
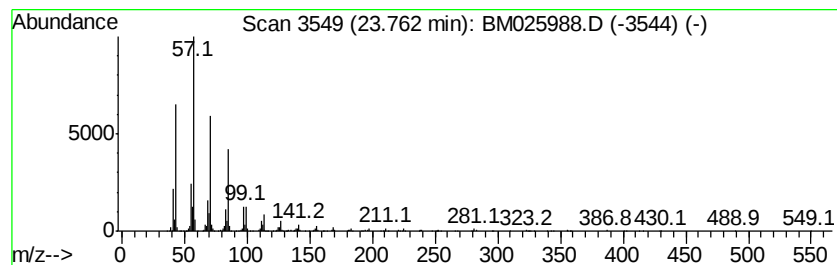
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.76	5.50 ng/ul	633750	Perylene-d12	23.22

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tricosane	324	C23H48	000638-67-5	94
2		Octadecane	254	C18H38	000593-45-3	91
3		Hexatriacontane	507	C36H74	000630-06-8	91
4		Octadecane	254	C18H38	000593-45-3	90
5		Heneicosane	296	C21H44	000629-94-7	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM051520\
Data File : BM025988.D
Acq On : 15 May 2020 14:45
Operator : MA/SJ
Sample : L2660-04
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
CB7C3

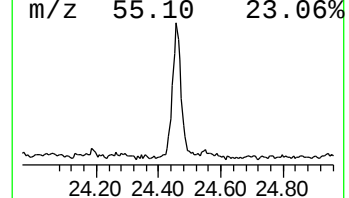
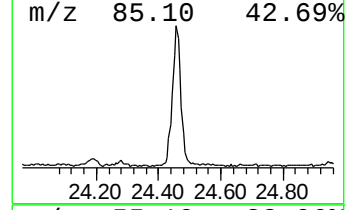
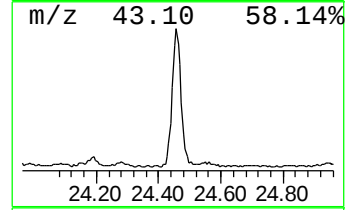
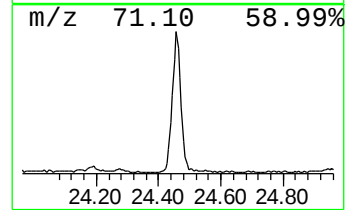
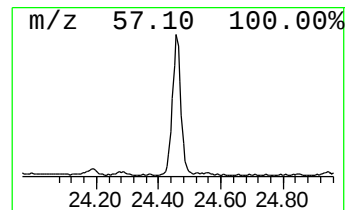
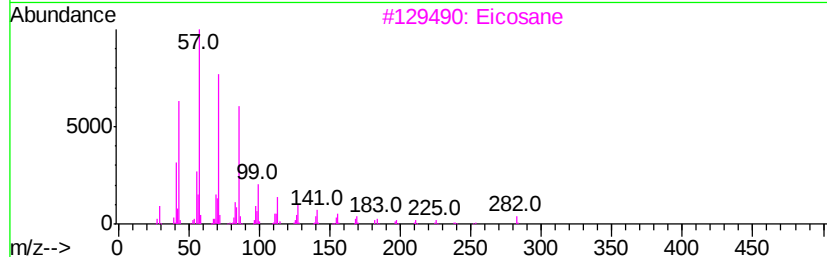
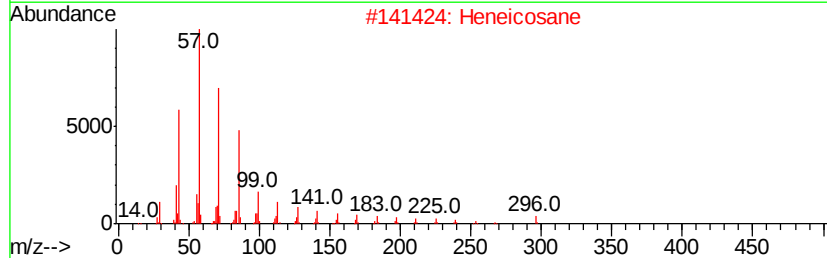
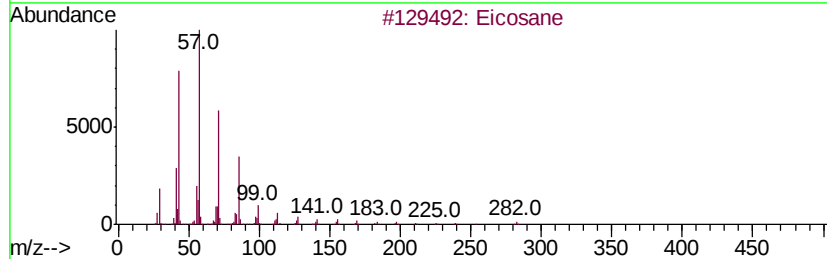
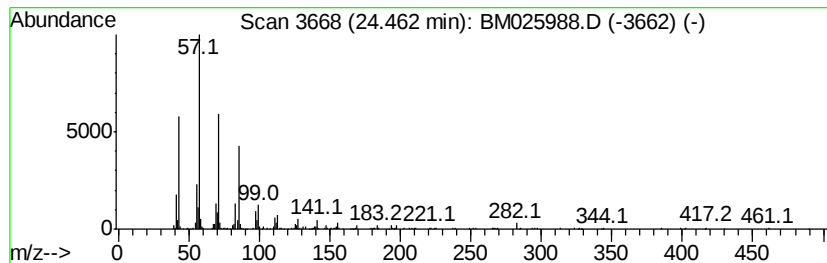
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.46	3.74 ng/ul	430856	Perylene-d12	23.22

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Eicosane	282	C20H42	000112-95-8	97
2			Heneicosane	296	C21H44	000629-94-7	95
3			Eicosane	282	C20H42	000112-95-8	94
4			Eicosane	282	C20H42	000112-95-8	94
5			Hexacosane	366	C26H54	000630-01-3	93



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM051520\
Data File : BM025988.D
Acq On : 15 May 2020 14:45
Operator : MA/SJ
Sample : L2660-04
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
CB7C3

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
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(DEL) Alkane: Str...	20.84	6.6	ng/ul	708356	5	21.10	2133020	20.0
(DEL) Alkane: Str...	21.29	2.8	ng/ul	299485	5	21.10	2133020	20.0
(DEL) Alkane: Str...	21.71	4.9	ng/ul	521398	5	21.10	2133020	20.0
(DEL) Alkane: Str...	22.15	6.2	ng/ul	656897	5	21.10	2133020	20.0
(DEL) Alkane: Str...	22.63	6.4	ng/ul	737547	6	23.22	2306490	20.0
(DEL) Alkane: Str...	23.16	6.8	ng/ul	788488	6	23.22	2306490	20.0
(DEL) Alkane: Str...	23.76	5.5	ng/ul	633750	6	23.22	2306490	20.0
(DEL) Alkane: Str...	24.46	3.7	ng/ul	430856	6	23.22	2306490	20.0