

Data Path : Z:\HPCHEM1\BNA M\DATA\BM102717\
 Data File : BM012481.D
 Acq On : 27 Oct 2017 11:56
 Operator : SJ/JU
 Sample : I5873-02
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 B-13(0-1) 101117

Integration Parameters: LSCINT.P

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

Filtering: 5
 Min Area: 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:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102417.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	3.869	129	133	142	rVB	152700	208370	1.24%	0.109%
2	4.093	166	171	180	rVB	843193	1126229	6.68%	0.588%
3	5.399	386	393	402	rVB	745020	1085367	6.43%	0.567%
4	5.528	409	415	428	rBV	3978813	7386430	43.79%	3.858%
5	5.899	466	478	489	rBV	1482103	2369213	14.05%	1.238%
6	7.040	664	672	688	rVB2	1681023	3293491	19.53%	1.720%
7	7.205	693	700	706	rBV	1477061	2227666	13.21%	1.164%
8	7.404	728	734	744	rVB	1118258	1785236	10.58%	0.933%
9	7.522	744	754	759	rBV2	81472	200260	1.19%	0.105%
10	7.875	804	814	820	rBV	2043523	3300523	19.57%	1.724%
11	8.110	847	854	860	rBV	729437	1153016	6.84%	0.602%
12	8.575	927	933	941	rBV	1995328	3101874	18.39%	1.620%
13	9.028	1003	1010	1018	rVB	1775897	2873496	17.04%	1.501%
14	9.746	1123	1132	1139	rBV	534357	953712	5.65%	0.498%
15	10.287	1215	1224	1232	rVB	668312	1144431	6.78%	0.598%
16	10.663	1279	1288	1294	rBV	2505256	4409404	26.14%	2.303%
17	10.798	1303	1311	1321	rBV	1331317	2391505	14.18%	1.249%
18	12.322	1564	1570	1579	rVB	107419	214283	1.27%	0.112%
19	13.334	1737	1742	1748	rVB2	140002	205047	1.22%	0.107%
20	13.404	1748	1754	1763	rBV	608041	943785	5.60%	0.493%
21	13.910	1831	1840	1844	rBV	3922953	5267275	31.23%	2.751%
22	13.957	1844	1848	1853	rVV	4035874	5282827	31.32%	2.759%
23	14.039	1857	1862	1868	rVB5	112847	204494	1.21%	0.107%
24	14.198	1883	1889	1898	rVB	4014013	6095746	36.14%	3.184%
25	14.404	1920	1924	1929	rBV	156965	223447	1.32%	0.117%
26	14.504	1932	1941	1947	rVV2	3340210	5287277	31.35%	2.762%
27	14.563	1947	1951	1960	rVB	171089	291546	1.73%	0.152%
28	14.898	2004	2008	2016	rVB2	204405	349125	2.07%	0.182%
29	15.304	2072	2077	2083	rBV	4839642	6332271	37.54%	3.308%
30	15.357	2083	2086	2091	rVB2	156410	198958	1.18%	0.104%
31	15.492	2103	2109	2115	rBV	5453996	7305322	43.31%	3.816%
32	15.545	2115	2118	2126	rVB2	199711	283940	1.68%	0.148%
33	15.628	2127	2132	2137	rBV	749962	912905	5.41%	0.477%
34	15.716	2143	2147	2152	rBV4	116353	243458	1.44%	0.127%

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35	15.875	2170	2174	2179	rVB	925484	1249389	7.41%	0.653%
36	16.222	2228	2233	2234	rBV2	268528	368144	2.18%	0.192%
37	16.239	2234	2236	2241	rVB	335039	425256	2.52%	0.222%
38	16.333	2248	2252	2256	rBV	246054	358651	2.13%	0.187%
39	16.386	2258	2261	2267	rVV2	239266	412853	2.45%	0.216%
40	16.451	2268	2272	2276	rVV2	212907	303891	1.80%	0.159%
41	16.598	2294	2297	2301	rVB3	191231	229932	1.36%	0.120%
42	16.651	2302	2306	2312	rVB4	275068	477912	2.83%	0.250%
43	16.716	2313	2317	2321	rBV	269940	362096	2.15%	0.189%
44	16.792	2326	2330	2333	rVB2	171951	203281	1.21%	0.106%
45	16.898	2344	2348	2356	rVB	211221	312755	1.85%	0.163%
46	17.010	2364	2367	2370	rVB	196396	224325	1.33%	0.117%
47	17.063	2372	2376	2385	rVB	355641	519098	3.08%	0.271%
48	17.245	2401	2407	2410	rBV2	3975895	5719291	33.91%	2.987%
49	17.286	2410	2414	2418	rVB	2444934	3334732	19.77%	1.742%
50	17.339	2418	2423	2428	rBV2	5068048	7352093	43.59%	3.840%
51	17.551	2455	2459	2466	rVB	1328445	1624351	9.63%	0.848%
52	17.645	2472	2475	2485	rVB	247106	435170	2.58%	0.227%
53	17.745	2489	2492	2499	rVB2	178616	255042	1.51%	0.133%
54	17.945	2523	2526	2534	rVB2	302297	454793	2.70%	0.238%
55	18.133	2552	2558	2560	rBV3	400531	813756	4.82%	0.425%
56	18.163	2561	2563	2567	rVB	437454	540810	3.21%	0.282%
57	18.310	2584	2588	2591	rBV2	331948	540597	3.20%	0.282%
58	18.616	2637	2640	2643	rBV	245122	296188	1.76%	0.155%
59	18.657	2644	2647	2657	rVB2	248487	415360	2.46%	0.217%
60	19.268	2747	2751	2752	rBV2	1270110	1333050	7.90%	0.696%
61	19.292	2752	2755	2762	rVB	3927362	5510567	32.67%	2.878%
62	19.627	2807	2812	2815	rBV	6426436	9383580	55.63%	4.901%
63	19.657	2815	2817	2823	rVB2	3451566	4667149	27.67%	2.438%
64	20.174	2903	2905	2910	rVB	729456	760541	4.51%	0.397%
65	20.551	2966	2969	2974	rVB	1381804	1553229	9.21%	0.811%
66	20.668	2987	2989	2998	rVB2	1110246	1473377	8.74%	0.770%
67	21.133	3064	3068	3073	rVB4	2208041	3002738	17.80%	1.568%
68	21.333	3098	3102	3106	rVB	16344786	16867486	100.00%	8.811%
69	21.409	3110	3115	3119	rVV2	5598179	9363896	55.51%	4.891%
70	21.451	3119	3122	3128	rVB	1761806	2242384	13.29%	1.171%
71	21.568	3139	3142	3148	rVB	2125037	2300932	13.64%	1.202%

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72	21.968	3206	3210	3215	rVV2	4219371	5105823	30.27%	2.667%
73	22.015	3215	3218	3226	rVB2	1905627	2589960	15.35%	1.353%
74	22.492	3297	3299	3305	rVB	1407079	1632087	9.68%	0.853%
75	23.021	3385	3389	3394	rBV2	2588816	4621290	27.40%	2.414%
76	23.574	3478	3483	3487	rBV2	4349215	7448564	44.16%	3.891%
77	23.721	3503	3508	3514	rVB	3170584	5707690	33.84%	2.981%

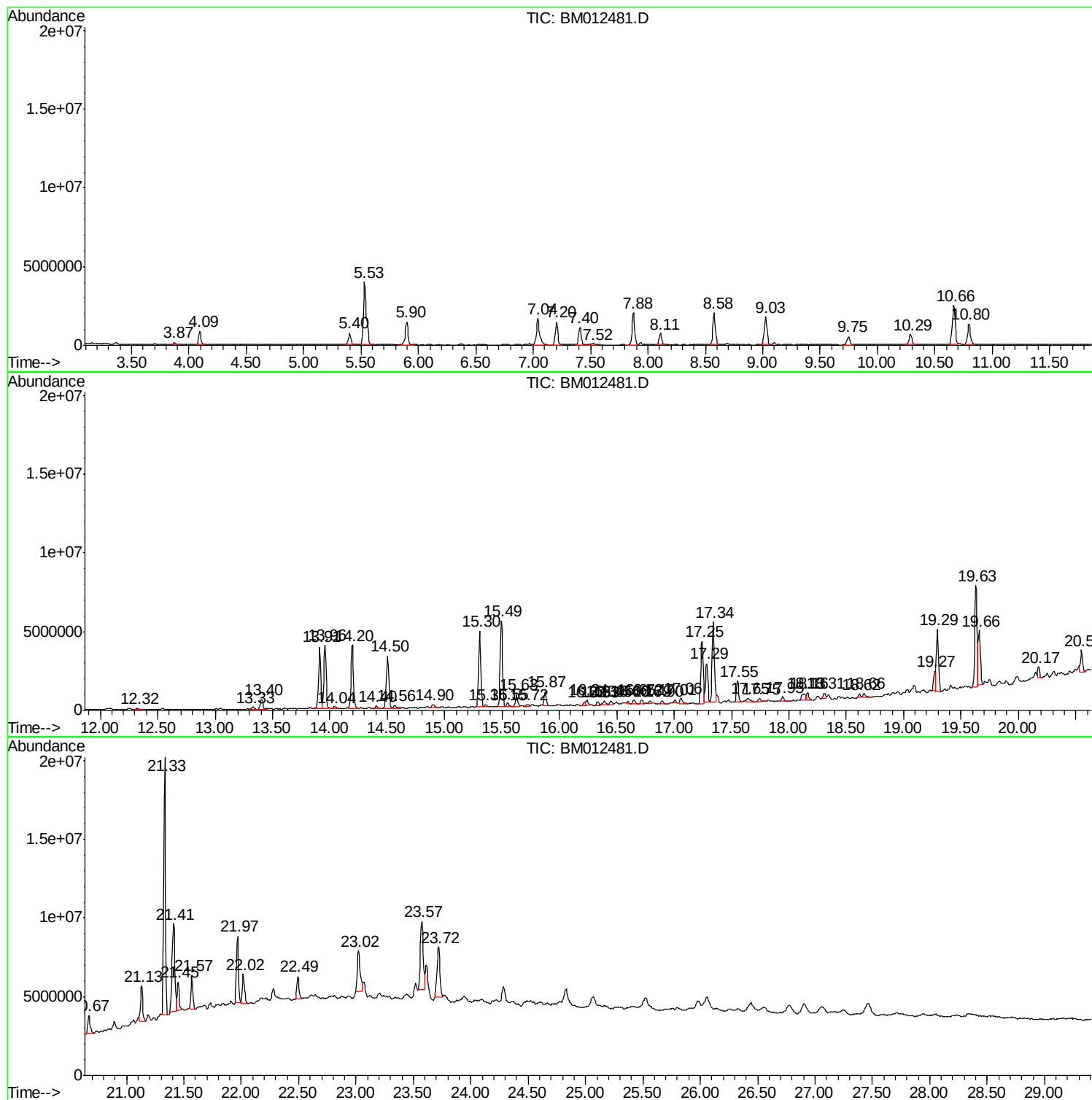
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Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM102417.M
Quant Title : SVOA CALIBRATION

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



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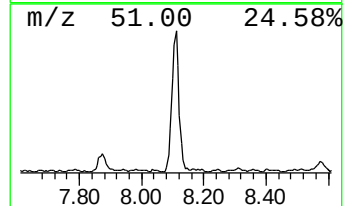
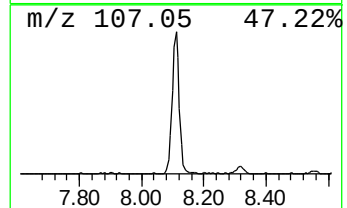
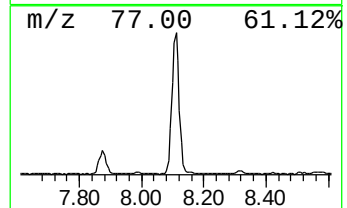
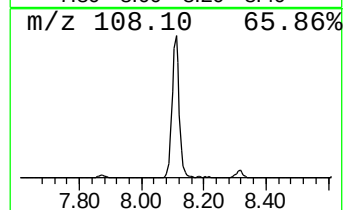
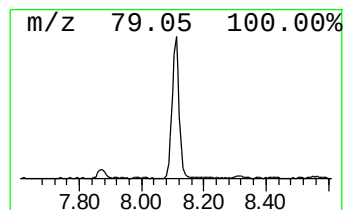
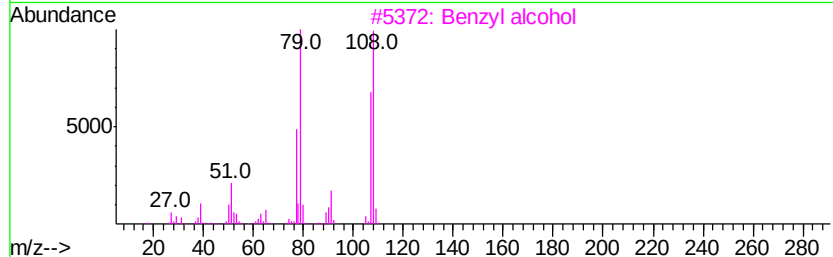
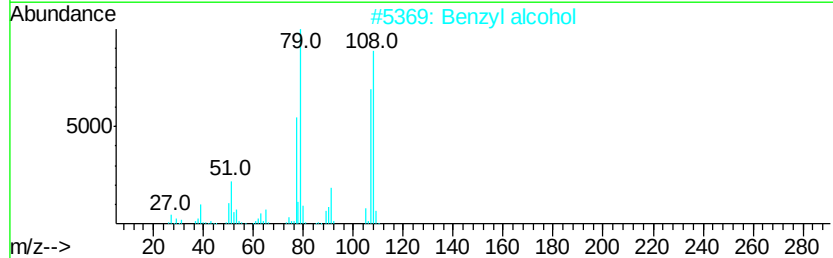
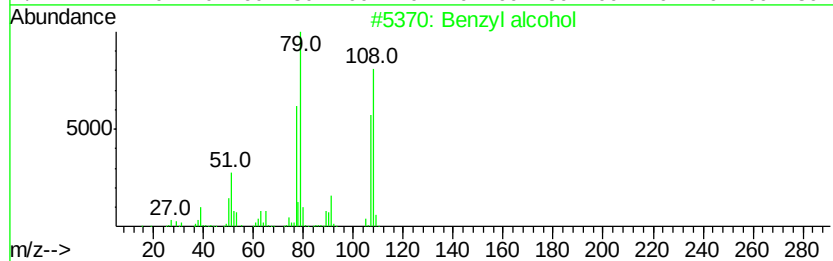
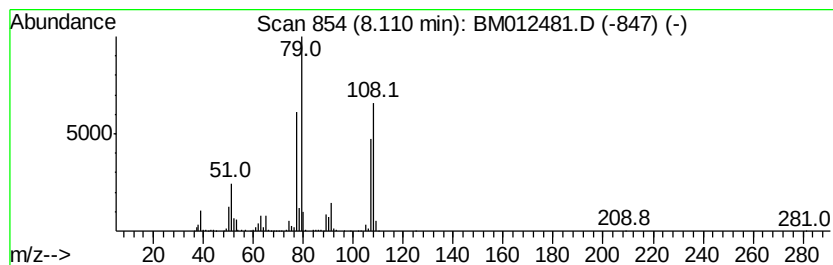
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Benzyl alcohol Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.11	6.99 ng/ul	1153020	1,4-Dichlorobenzene-d4	7.88

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzyl alcohol	108	C7H8O	000100-51-6	98
2		Benzyl alcohol	108	C7H8O	000100-51-6	97
3		Benzyl alcohol	108	C7H8O	000100-51-6	96
4		Benzyl alcohol	108	C7H8O	000100-51-6	95
5		N-Cbz-glycylglycine p-nitropheny...	387	C18H17N3O7	013574-81-7	91



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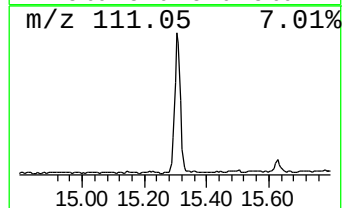
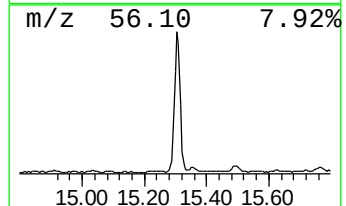
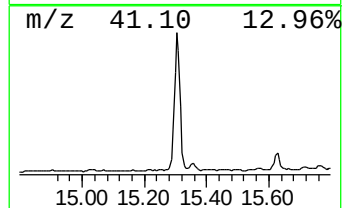
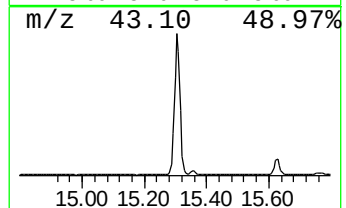
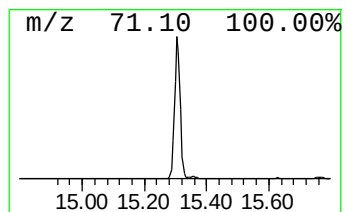
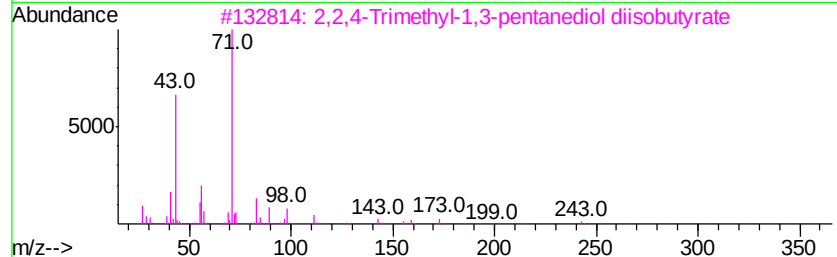
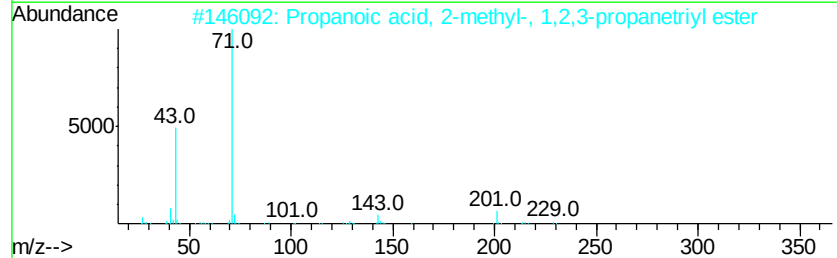
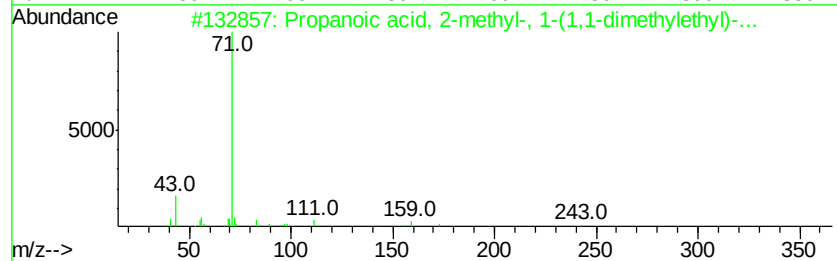
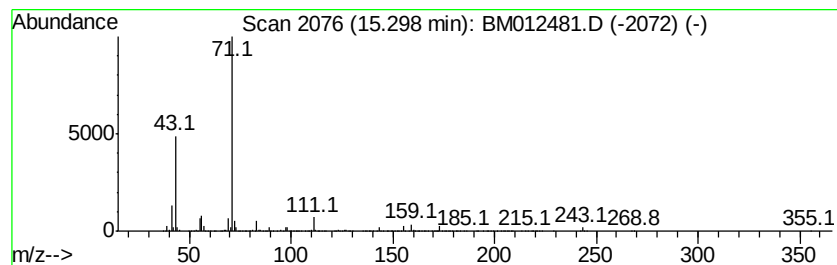
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 Propanoic acid, 2-methyl-, ... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.30	23.95 ng/ul	6332270	Acenaphthene-d10	14.50
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Propanoic acid, 2-methyl-, 1-(1,...	286 C16H30O4	074381-40-1	64
2	Propanoic acid, 2-methyl-, 1,2,3...	302 C15H26O6	014295-64-8	59
3	2,2,4-Trimethyl-1,3-pentanediol ...	286 C16H30O4	006846-50-0	56
4	Diethylmalonic acid, monochlorid...	262 C12H19ClO4	1000370-65-0	45
5	Butanoic acid, 2-butoxy-1-methyl...	216 C11H20O4	007492-70-8	45



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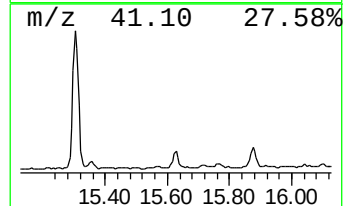
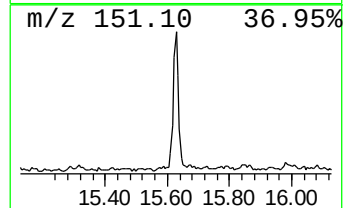
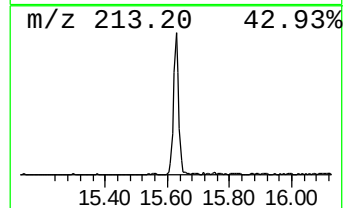
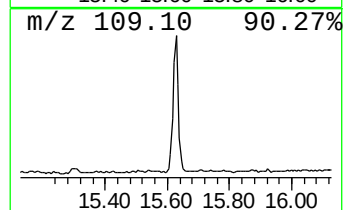
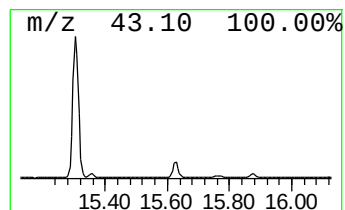
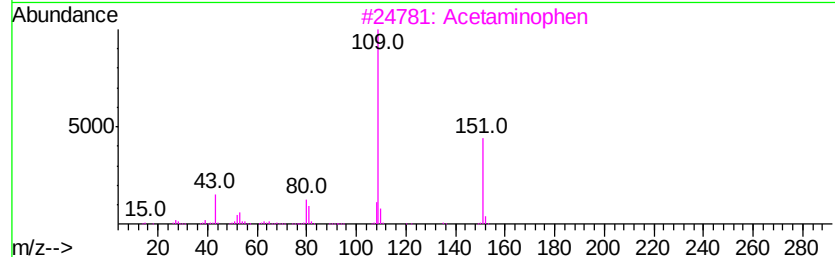
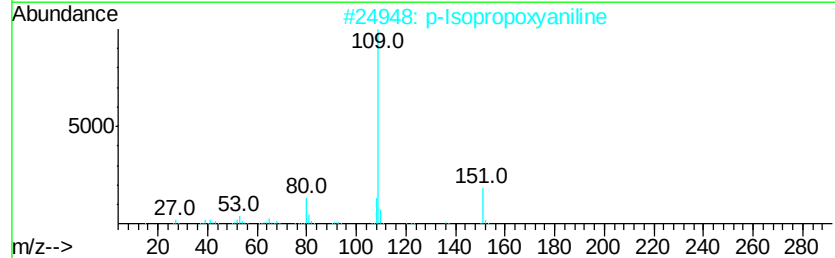
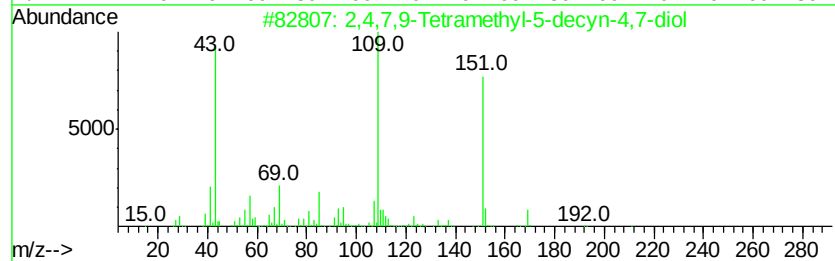
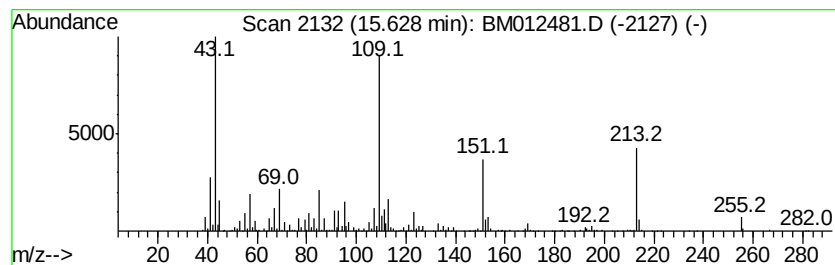
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 2,4,7,9-Tetramethyl-5-decyn... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.63	3.45 ng/ul	912905	Acenaphthene-d10	14.50

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,4,7,9-Tetramethyl-5-decyn-4,7-...	226	C14H26O2	000126-86-3	58		
2	p-Isopropoxyaniline	151	C9H13NO	007664-66-6	38		
3	Acetaminophen	151	C8H9NO2	000103-90-2	38		
4	Metacetamol	151	C8H9NO2	000621-42-1	38		
5	Acetaminophen	151	C8H9NO2	000103-90-2	35		



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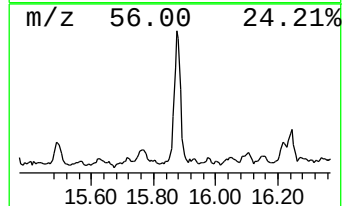
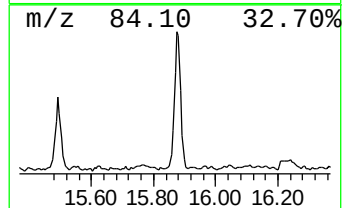
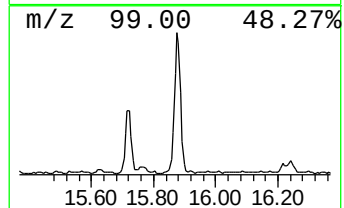
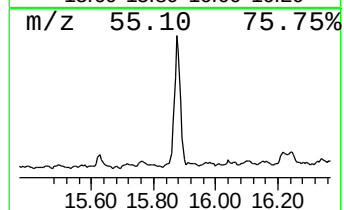
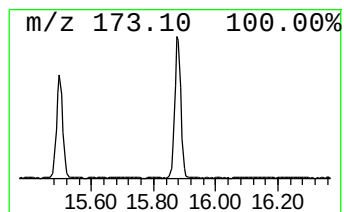
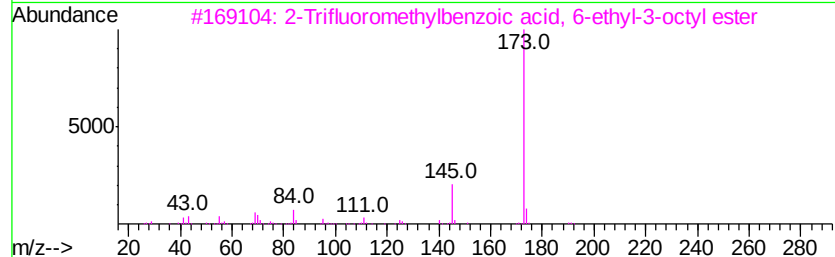
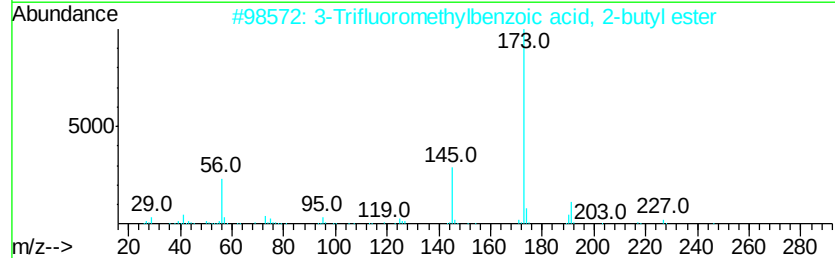
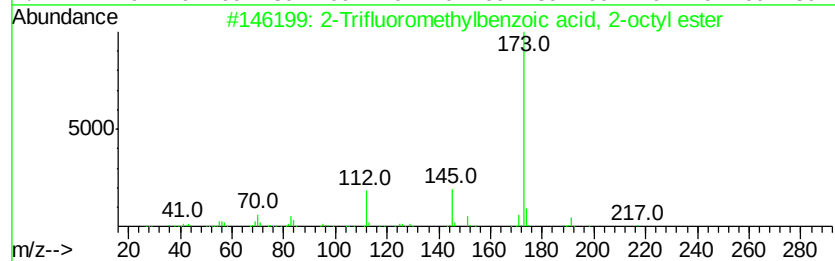
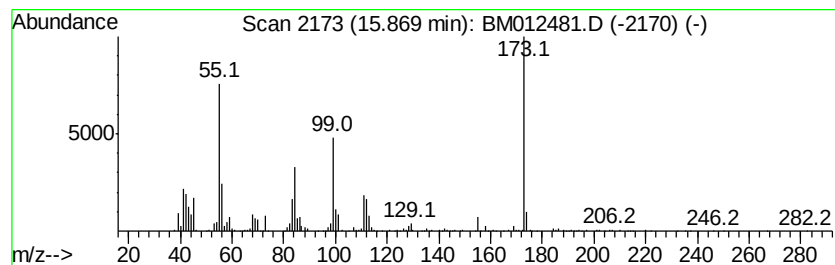
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 unknown-01 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.87	4.73 ng/ul	1249390	Acenaphthene-d10	14.50

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Trifluoromethylbenzoic acid, 2...	302	C16H21F3O2	1000282-64-8	35
2		3-Trifluoromethylbenzoic acid, 2...	246	C12H13F3O2	1000282-68-1	35
3		2-Trifluoromethylbenzoic acid, 6...	330	C18H25F3O2	1000282-65-0	25
4		2,6-Difluoro-3-methylbenzoic aci...	368	C22H34F2O2	1000338-85-1	23
5		1,2-Benzenediol, 0-(6-chlorohexa...	414	C20H18ClF3O4	1000329-80-9	16



Data Path : Z:\HPCHEM1\BNA M\DATA\BM102717\
Data File : BM012481.D
Acq On : 27 Oct 2017 11:56
Operator : SJ/JU
Sample : I5873-02
Misc :
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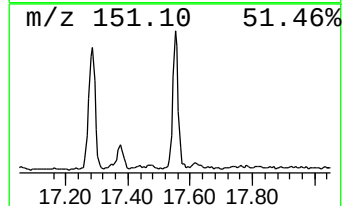
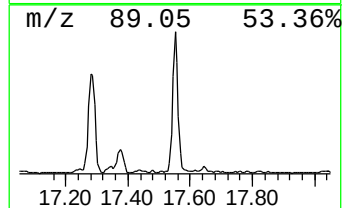
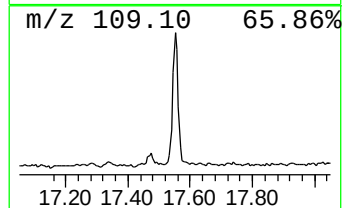
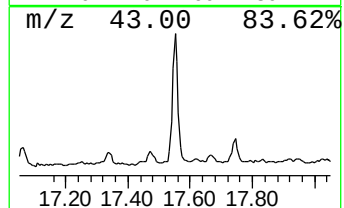
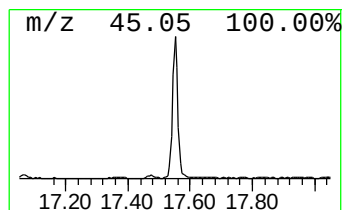
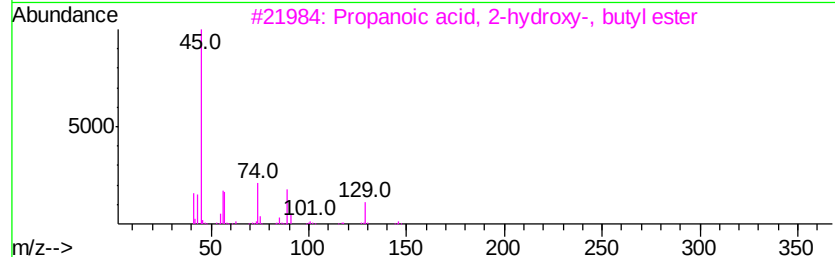
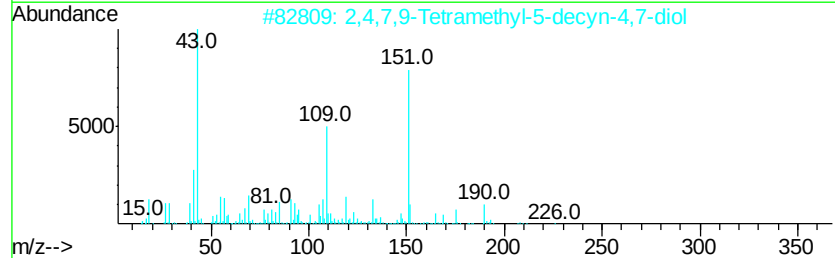
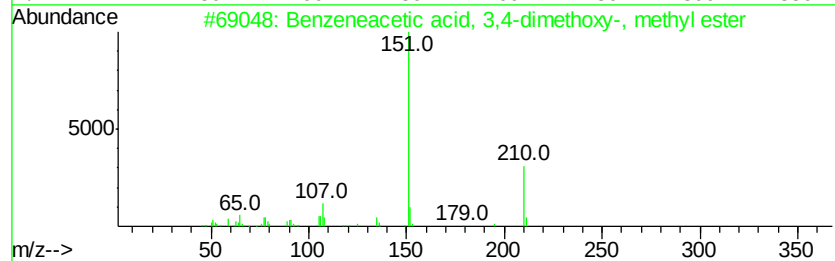
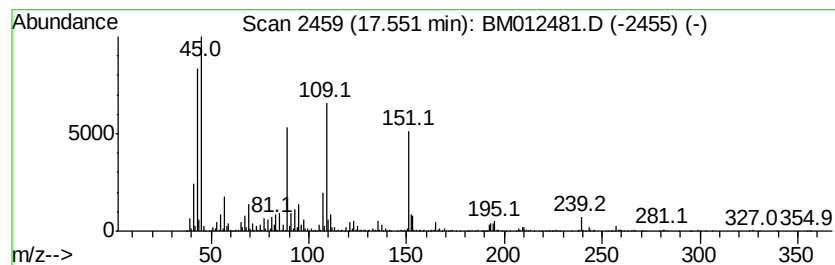
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 unknown-02 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.		
17.55	5.68 ng/ul	1624350	Phenanthrene-d10	17.25		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzeneacetic acid, 3,4-dimethox...	210	C11H14O4	015964-79-1	30
2		2,4,7,9-Tetramethyl-5-decyn-4,7-...	226	C14H26O2	000126-86-3	27
3		Propanoic acid, 2-hydroxy-, buty...	146	C7H14O3	000138-22-7	22
4		Propanoic acid, 2-hydroxy-, buty...	146	C7H14O3	000138-22-7	22
5		5-Methoxymethoxyhexa-2,3-diene	142	C8H14O2	1000190-45-1	22



Data Path : Z:\HPCHEM1\BNA M\DATA\BM102717\
Data File : BM012481.D
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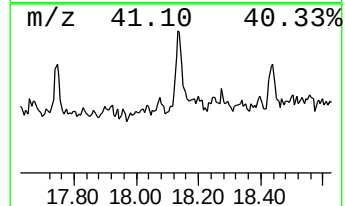
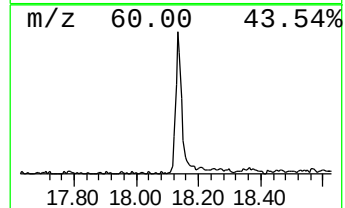
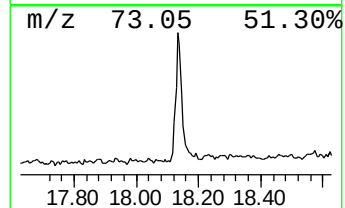
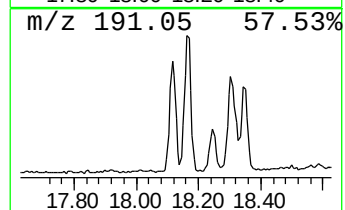
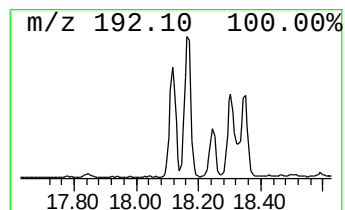
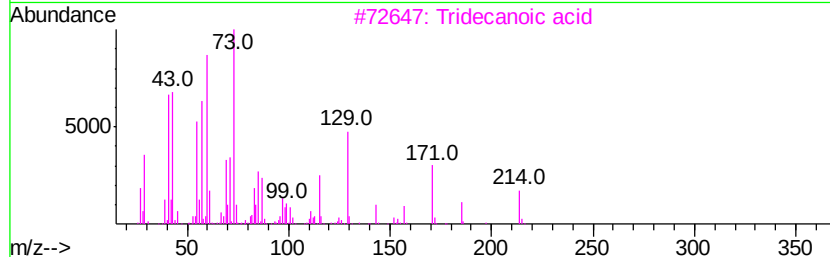
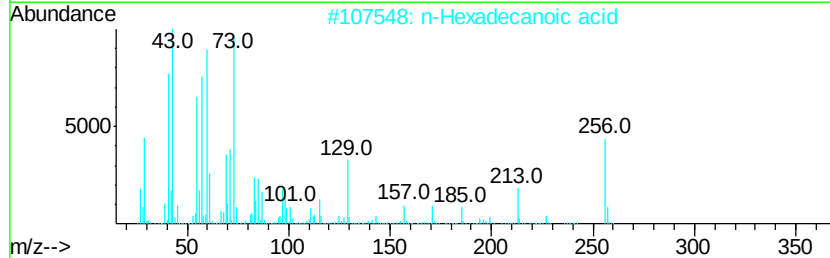
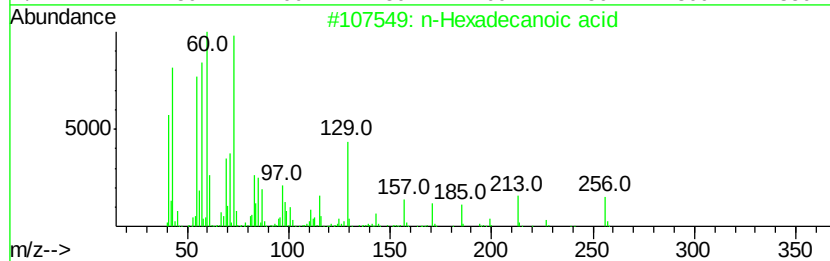
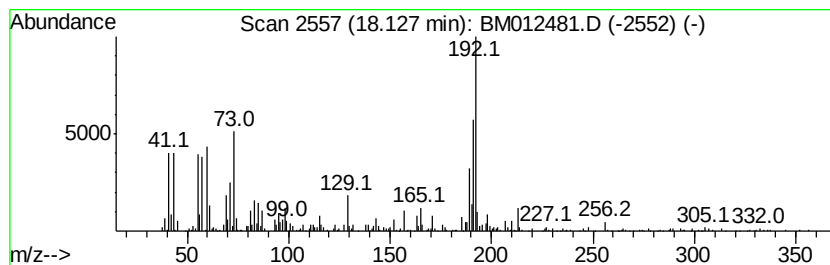
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 n-Hexadecanoic acid Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.13	2.85 ng/ul	813756	Phenanthrene-d10	17.25

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
3		Tridecanoic acid	214	C13H26O2	000638-53-9	95
4		Tridecanoic acid	214	C13H26O2	000638-53-9	89
5		Pentadecanoic acid	242	C15H30O2	001002-84-2	70



Data Path : Z:\HPCHEM1\BNA M\DATA\BM102717\
Data File : BM012481.D
Acq On : 27 Oct 2017 11:56
Operator : SJ/JU
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Misc :
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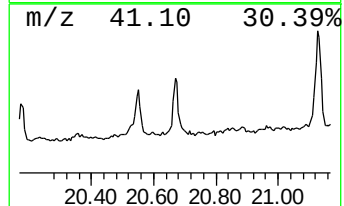
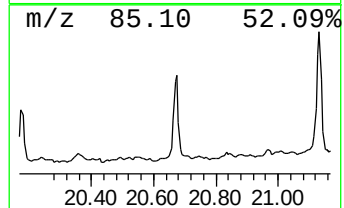
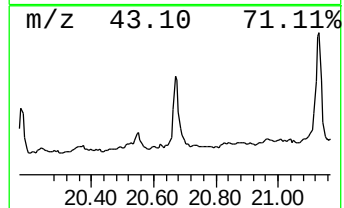
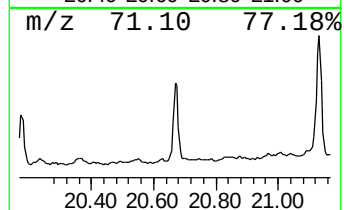
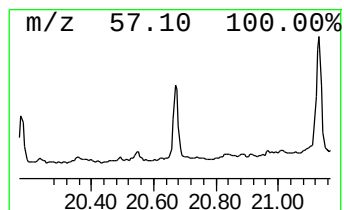
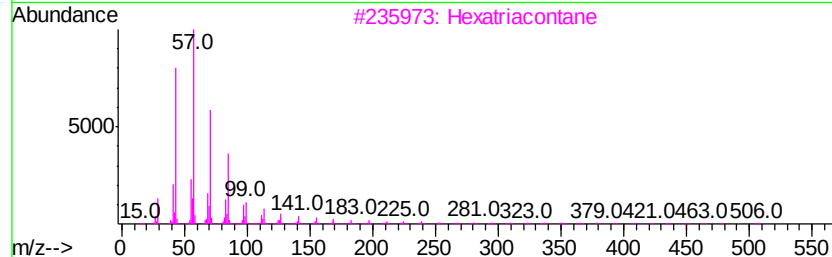
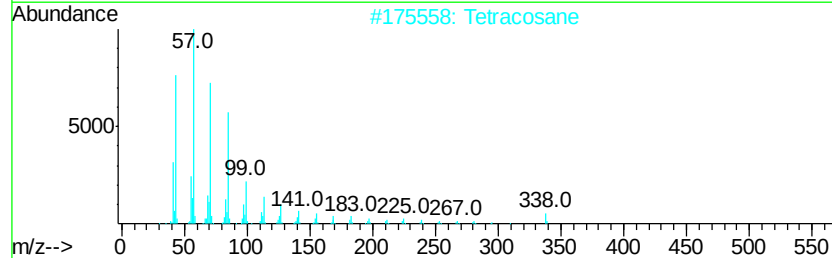
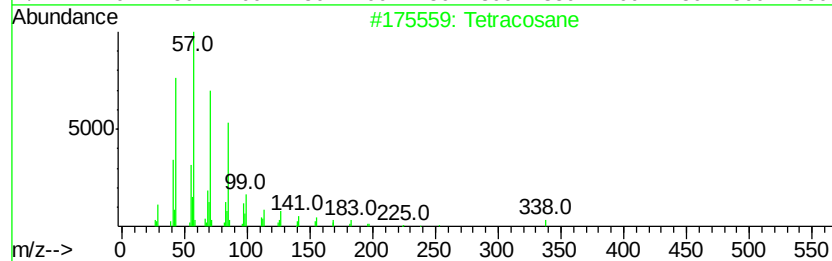
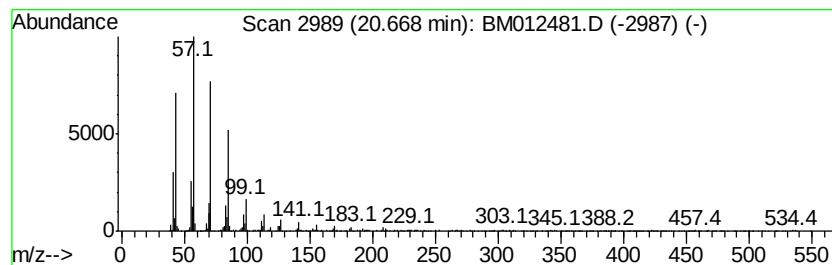
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Quant Title : SVOA CALIBRATION

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

Peak Number 12 (DEL) Alkane: Straight-Chai... Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.67	3.15 ng/ul	1473380	Chrysene-d12	21.42

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetracosane	338	C24H50	000646-31-1	97
2			Tetracosane	338	C24H50	000646-31-1	95
3			Hexatriacontane	507	C36H74	000630-06-8	95
4			Tetracosane	338	C24H50	000646-31-1	93
5			Tetracosane	338	C24H50	000646-31-1	90



Data Path : Z:\HPCHEM1\BNA M\DATA\BM102717\
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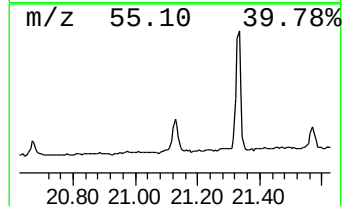
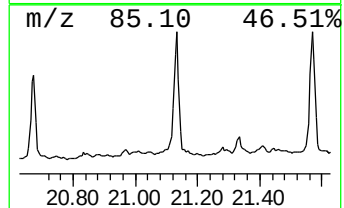
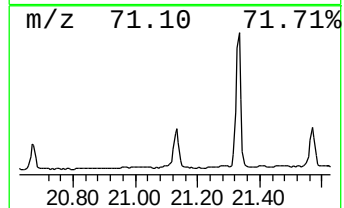
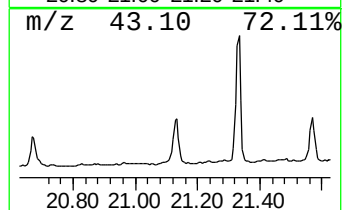
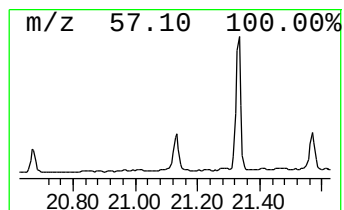
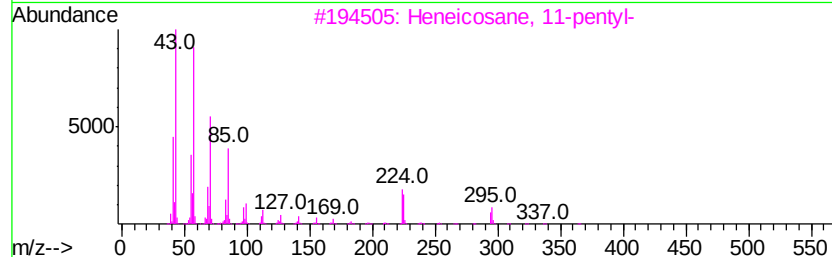
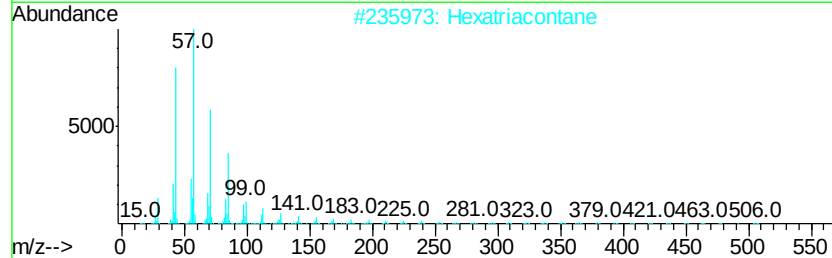
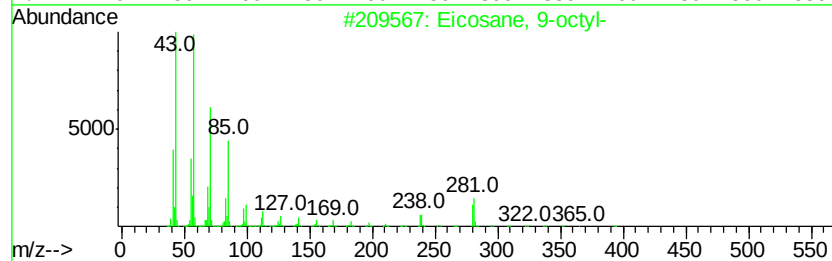
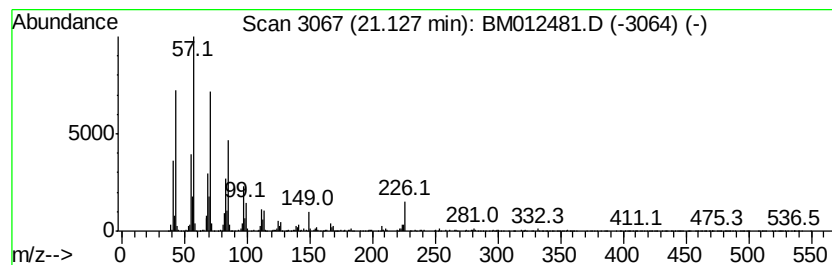
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Quant Title : SVOA CALIBRATION

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

Peak Number 13 (DEL) Alkane: Straight-Chai... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD		R.T.
21.13	6.41 ng/ul	3002740	Chrysene-d12		21.42
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane, 9-octyl-	394	C28H58	013475-77-9	91
2	Hexatriacontane	507	C36H74	000630-06-8	91
3	Heneicosane, 11-pentyl-	366	C26H54	014739-72-1	87
4	Hexadecane	226	C16H34	000544-76-3	87
5	Docosane, 11-butyl-	366	C26H54	013475-76-8	87



Data Path : Z:\HPCHEM1\BNA M\DATA\BM102717\
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Acq On : 27 Oct 2017 11:56
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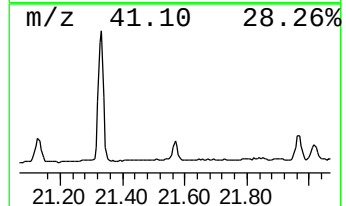
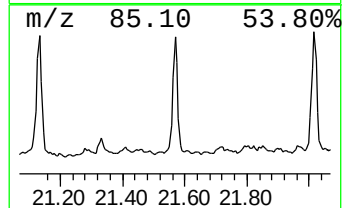
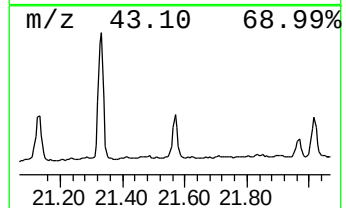
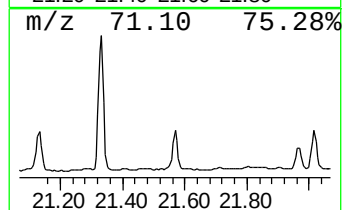
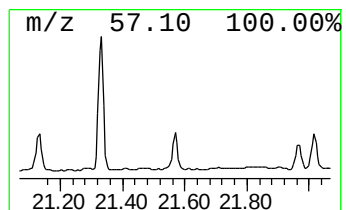
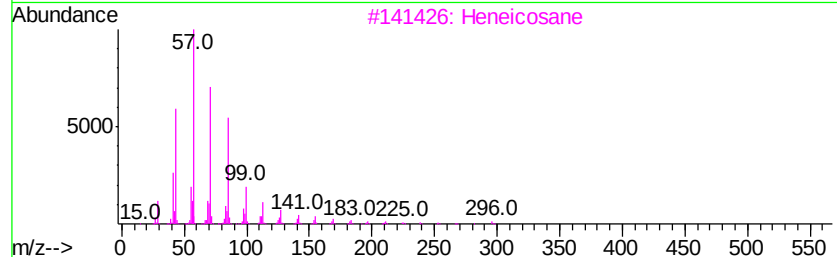
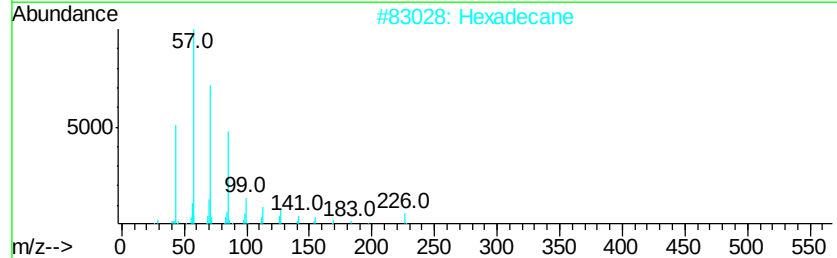
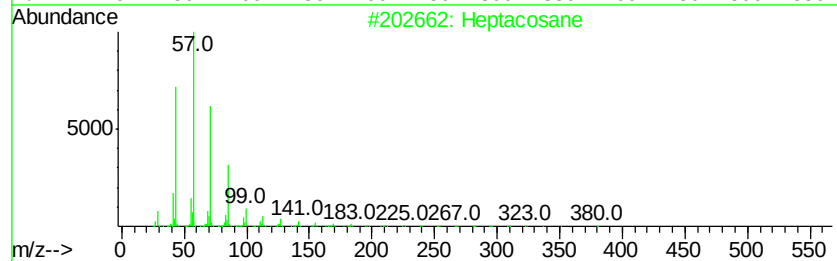
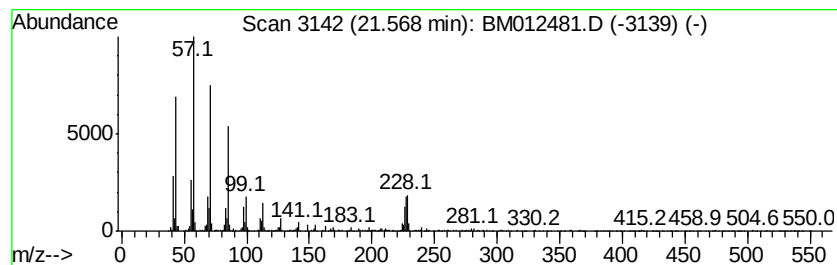
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Quant Title : SVOA CALIBRATION

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

Peak Number 14 (DEL) Alkane: Straight-Chai... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.57	4.91 ng/ul	2300930	Chrysene-d12	21.42

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptacosane	380	C27H56	000593-49-7	91
2			Hexadecane	226	C16H34	000544-76-3	81
3			Heneicosane	296	C21H44	000629-94-7	76
4			Hexacosane	366	C26H54	000630-01-3	76
5			Octadecane, 2-methyl-	268	C19H40	001560-88-9	74



Data Path : Z:\HPCHEM1\BNA M\DATA\BM102717\
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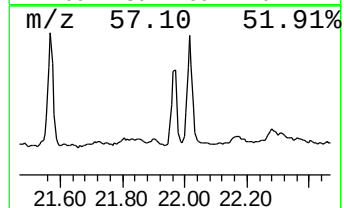
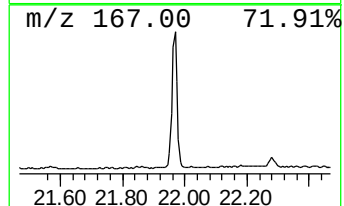
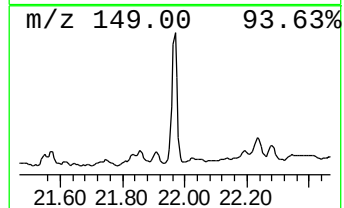
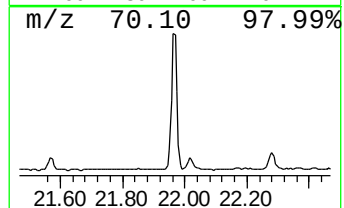
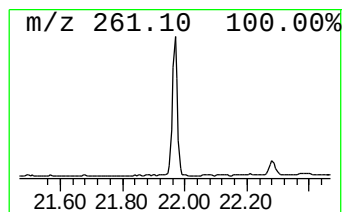
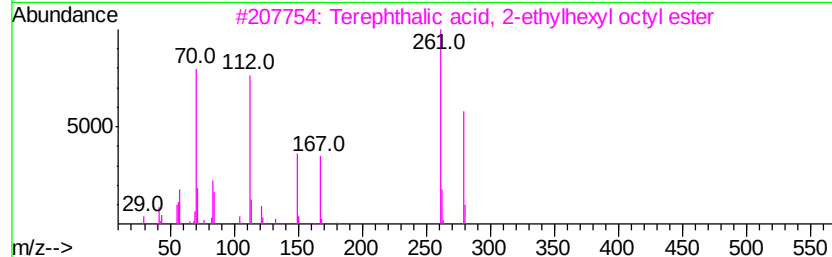
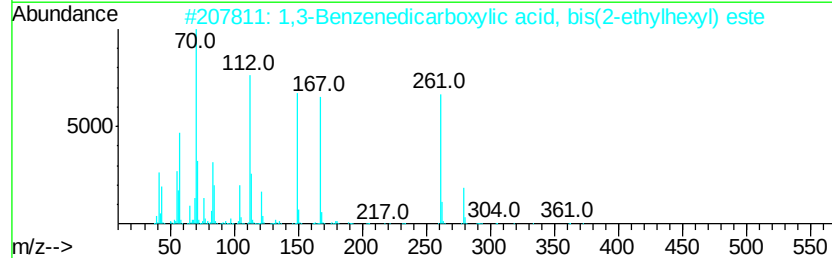
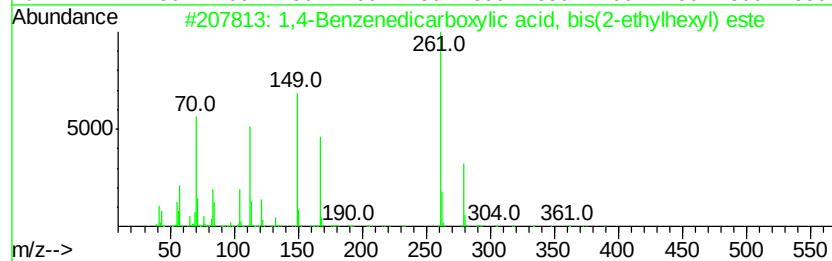
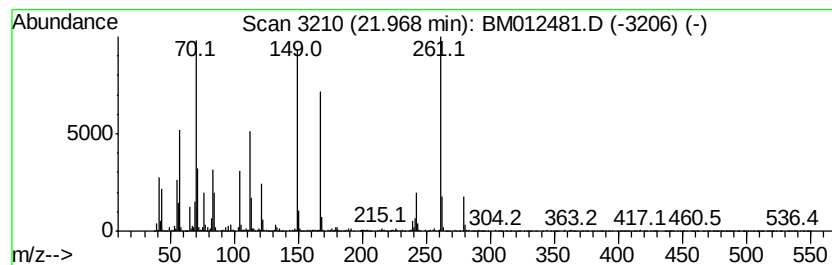
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Quant Title : SVOA CALIBRATION

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

Peak Number 15 1,4-Benzenedicarboxylic aci... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.97	10.91 ng/ul	5105820	Chrysene-d12	21.42

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,4-Benzenedicarboxylic acid, bi...	390	C24H38O4	006422-86-2	64
2		1,3-Benzenedicarboxylic acid, bi...	390	C24H38O4	000137-89-3	59
3		Terephthalic acid, 2-ethylhexyl ...	390	C24H38O4	1000324-00-5	43
4		Isophthalic acid, 2-chloro-5-met...	402	C23H27ClO4	1000356-57-0	35
5		Isophthalic acid, 4-chlorophenyl...	388	C22H25ClO4	1000344-58-3	35



Data Path : Z:\HPCHEM1\BNA M\DATA\BM102717\
Data File : BM012481.D
Acq On : 27 Oct 2017 11:56
Operator : SJ/JU
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ALS Vial : 6 Sample Multiplier: 1

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ClientSampleId :
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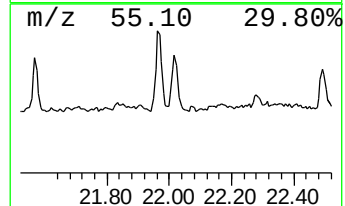
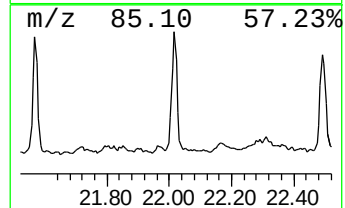
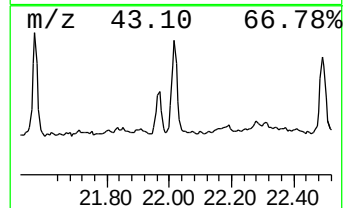
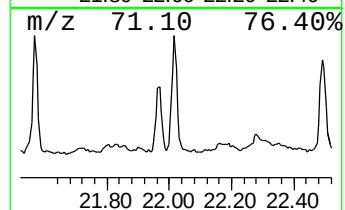
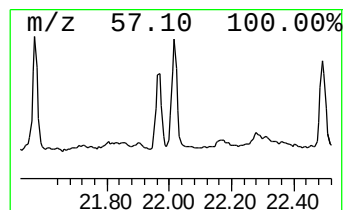
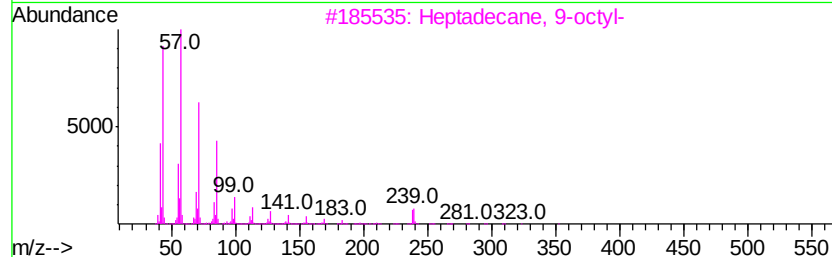
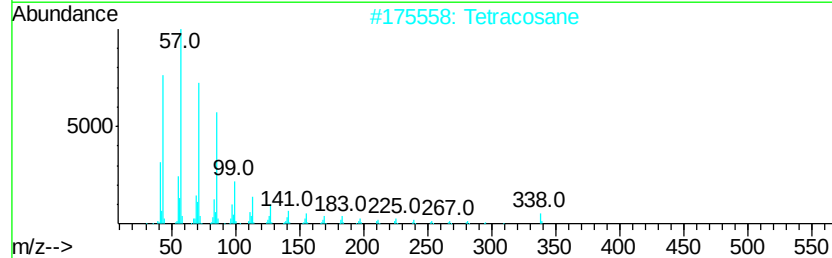
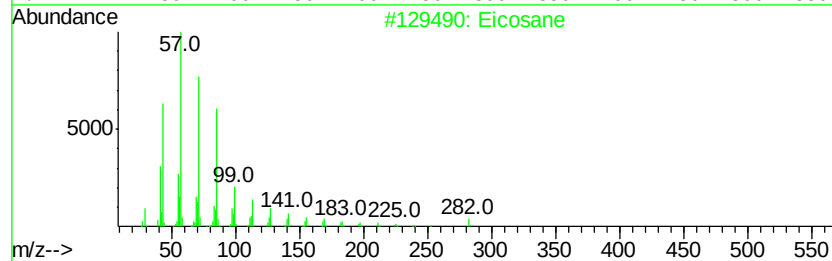
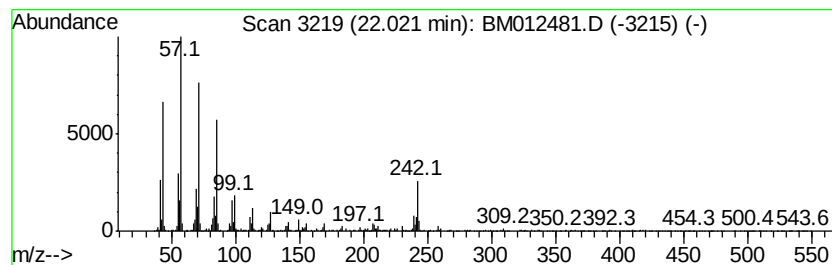
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Quant Title : SVOA CALIBRATION

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

Peak Number 16 (DEL) Alkane: Straight-Chai... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.02	5.53 ng/ul	2589960	Chrysene-d12	21.42

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Eicosane	282	C20H42	000112-95-8	97
2			Tetracosane	338	C24H50	000646-31-1	97
3			Heptadecane, 9-octyl-	352	C25H52	007225-64-1	95
4			Tridecane, 1-iodo-	310	C13H27I	035599-77-0	93
5			Octacosane	394	C28H58	000630-02-4	91



Data Path : Z:\HPCHEM1\BNA M\DATA\BM102717\
Data File : BM012481.D
Acq On : 27 Oct 2017 11:56
Operator : SJ/JU
Sample : I5873-02
Misc :
ALS Vial : 6 Sample Multiplier: 1

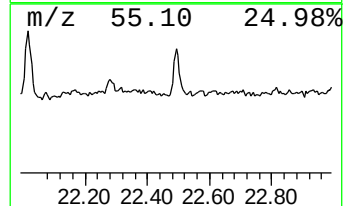
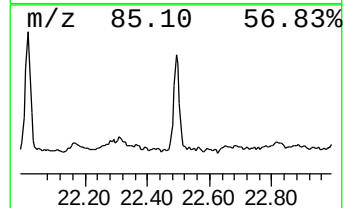
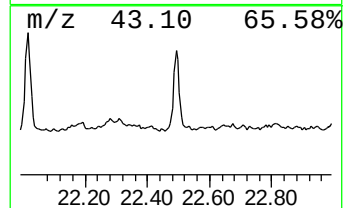
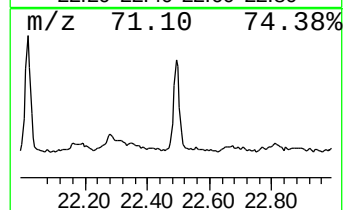
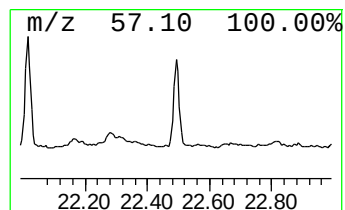
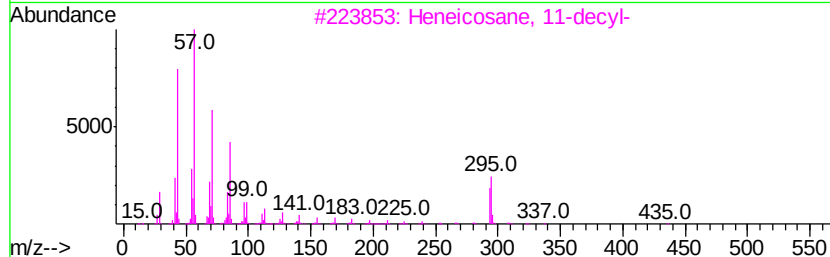
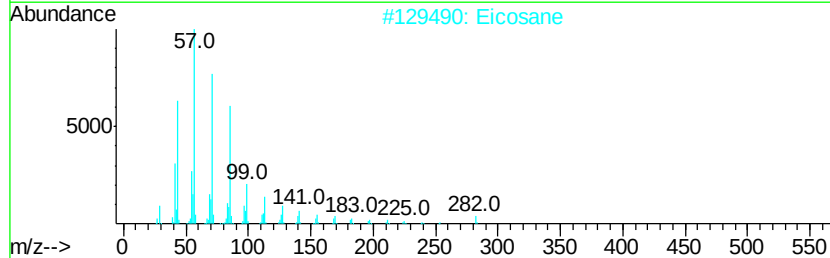
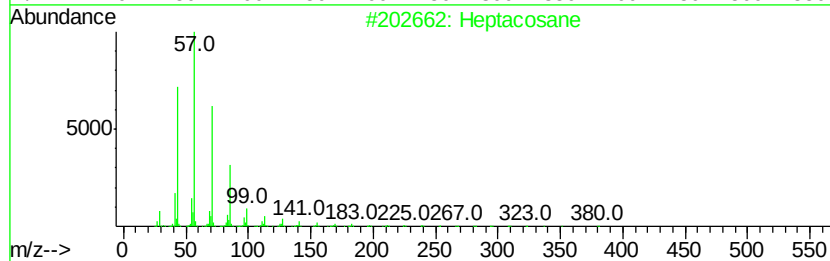
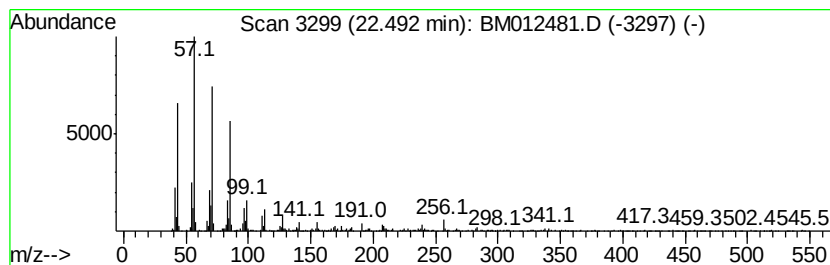
Instrument :
BNA_M
ClientSampleId :
B-13(0-1) 101117

Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM102417.M
Quant Title : SVOA CALIBRATION

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

Peak Number 17 (DEL) Alkane: Straight-Chai... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD		R.T.
22.49	3.49 ng/ul	1632090	Chrysene-d12		21.42
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptacosane	380	C27H56	000593-49-7	94
2	Eicosane	282	C20H42	000112-95-8	93
3	Heneicosane, 11-decyl-	437	C31H64	055320-06-4	93
4	Octadecane, 2-methyl-	268	C19H40	001560-88-9	93
5	Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	93



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM102717\
 Data File : BM012481.D
 Acq On : 27 Oct 2017 11:56
 Operator : SJ/JU
 Sample : I5873-02
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 B-13(0-1)_101117

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102417.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
Benzyl alcohol	8.11	7.0	ng/ul	1153020	1	7.88	3300520	20.0
Propanoic acid, 2...	15.30	23.9	ng/ul	6332270	3	14.50	5287280	20.0
2,4,7,9-Tetrameth...	15.63	3.5	ng/ul	912905	3	14.50	5287280	20.0
unknown-01	15.87	4.7	ng/ul	1249390	3	14.50	5287280	20.0
unknown-02	17.55	5.7	ng/ul	1624350	4	17.25	5719290	20.0
n-Hexadecanoic acid	18.13	2.9	ng/ul	813756	4	17.25	5719290	20.0
(DEL) Alkane: Str...	20.67	3.1	ng/ul	1473380	5	21.42	9363900	20.0
(DEL) Alkane: Str...	21.13	6.4	ng/ul	3002740	5	21.42	9363900	20.0
(DEL) Alkane: Str...	21.57	4.9	ng/ul	2300930	5	21.42	9363900	20.0
1,4-Benzenedicarb...	21.97	10.9	ng/ul	5105820	5	21.42	9363900	20.0
(DEL) Alkane: Str...	22.02	5.5	ng/ul	2589960	5	21.42	9363900	20.0
(DEL) Alkane: Str...	22.49	3.5	ng/ul	1632090	5	21.42	9363900	20.0