

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF080218\
 Data File : BF107924.D
 Acq On : 2 Aug 2018 22:01
 Operator : JU/SJ
 Sample : J4285-21
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 KG-PIT-6

Integration Parameters: rteint.p

Integrator: RTE

Smoothing : ON

Sampling : 1

Start Thrs: 0.2

Stop Thrs : 0

Filtering: 5

Min Area: 3 % 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_F\METHODS\8270-BF073018.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.190	481	485	495	rBV	274358	393680	12.86%	1.542%
2	5.596	551	554	572	rBV	896867	1952916	63.79%	7.651%
3	6.601	721	725	743	rBV	993092	2278635	74.43%	8.927%
4	6.731	743	747	781	rVV	1739547	3061579	100.00%	11.994%
5	6.954	782	785	807	rVV	325531	793134	25.91%	3.107%
6	7.107	807	811	831	rVV	1583498	1985211	64.84%	7.777%
7	7.525	878	882	904	rBV	596409	1500981	49.03%	5.880%
8	8.237	999	1003	1021	rBV	555332	950084	31.03%	3.722%
9	9.307	1181	1185	1194	rBV	1865764	2272928	74.24%	8.904%
10	9.701	1246	1252	1258	rBV	123592	170998	5.59%	0.670%
11	9.895	1281	1285	1289	rVB2	35655	38142	1.25%	0.149%
12	9.989	1298	1301	1322	rBV	674234	873649	28.54%	3.423%
13	10.395	1365	1370	1373	rBV2	54690	57218	1.87%	0.224%
14	10.619	1403	1408	1410	rBV	27191	34688	1.13%	0.136%
15	10.789	1433	1437	1448	rBV	781317	1213392	39.63%	4.754%
16	10.872	1448	1451	1457	rVB2	68307	108512	3.54%	0.425%
17	11.319	1524	1527	1529	rBV	52687	46364	1.51%	0.182%
18	11.348	1529	1532	1534	rVB	40132	36298	1.19%	0.142%
19	11.489	1552	1556	1567	rBV	494256	842593	27.52%	3.301%
20	11.748	1597	1600	1603	rBV	59724	45959	1.50%	0.180%
21	11.995	1638	1642	1646	rBV5	59021	119710	3.91%	0.469%
22	12.095	1657	1659	1663	rBV3	25632	31575	1.03%	0.124%
23	12.154	1667	1669	1672	rBV	51556	46530	1.52%	0.182%
24	12.436	1714	1717	1720	rBV	44018	45523	1.49%	0.178%
25	12.542	1732	1735	1738	rBV	83072	92010	3.01%	0.360%
26	12.683	1757	1759	1761	rBV	43312	47637	1.56%	0.187%
27	12.707	1761	1763	1769	rVB	101205	140754	4.60%	0.551%
28	12.919	1797	1799	1800	rBV2	50036	42484	1.39%	0.166%
29	12.942	1800	1803	1807	rVB	107571	129912	4.24%	0.509%
30	13.072	1821	1825	1831	rBV	1582685	1712077	55.92%	6.707%
31	13.272	1857	1859	1862	rBV2	77256	66149	2.16%	0.259%
32	13.395	1878	1880	1884	rBV2	107476	129177	4.22%	0.506%
33	13.589	1909	1913	1915	rBV	708727	696836	22.76%	2.730%
34	13.730	1935	1937	1944	rVB2	155614	226376	7.39%	0.887%

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Integration Parameters: rteint.p
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % 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

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Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	14.054	1989	1992	1996	rBV	310242	480337	15.69%	1.882%
36	14.142	2004	2007	2014	rBV	673655	1035726	33.83%	4.058%
37	14.366	2042	2045	2051	rBV2	356059	563161	18.39%	2.206%
38	14.666	2093	2096	2104	rVB	357460	567656	18.54%	2.224%
39	15.677	2265	2268	2278	rVB	378982	695476	22.72%	2.725%

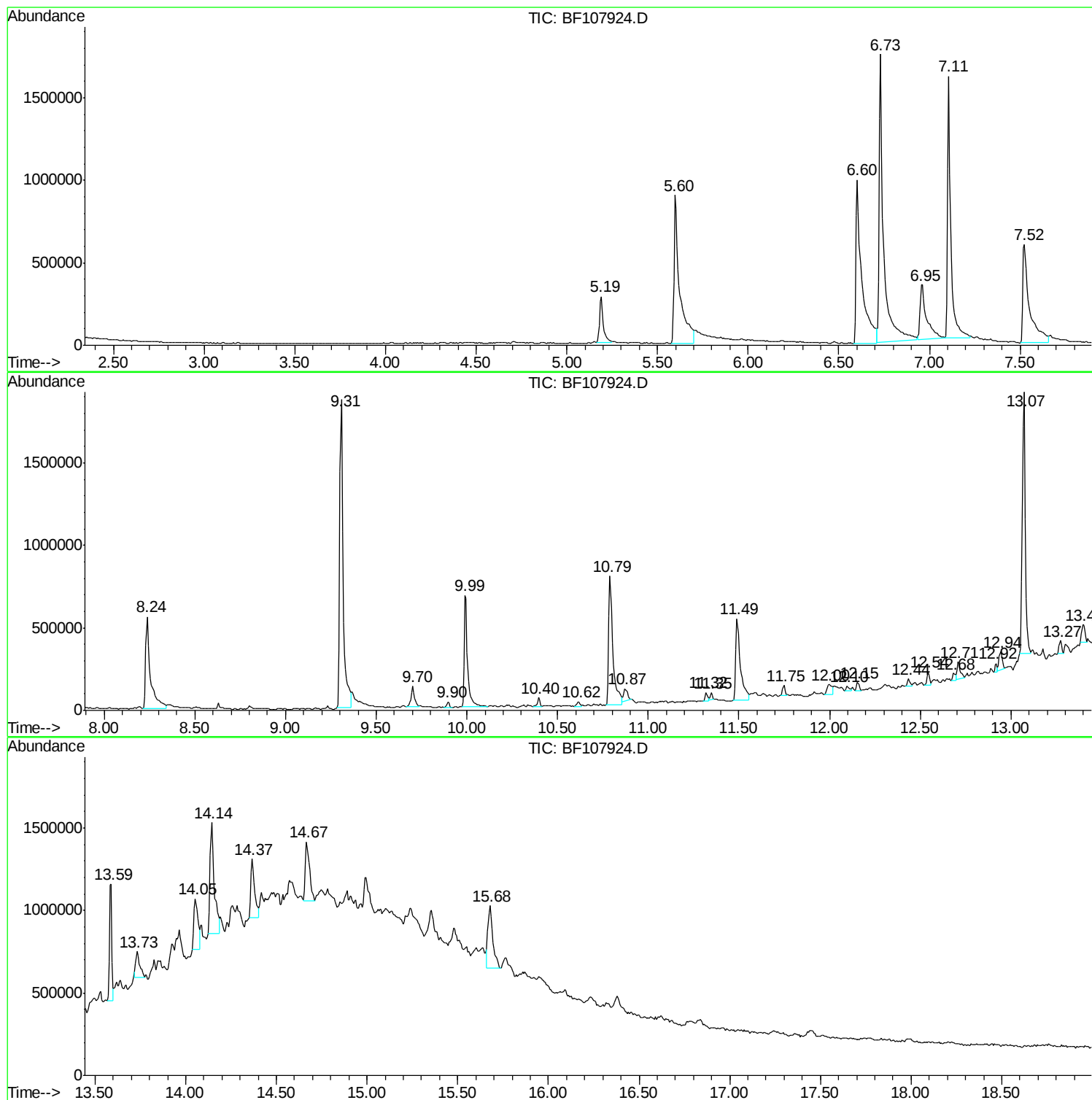
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF073018.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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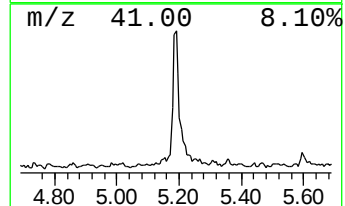
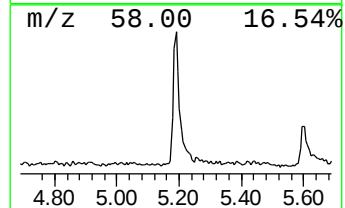
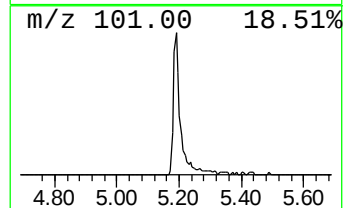
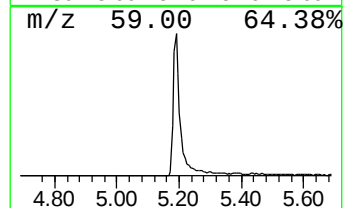
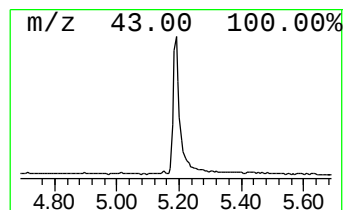
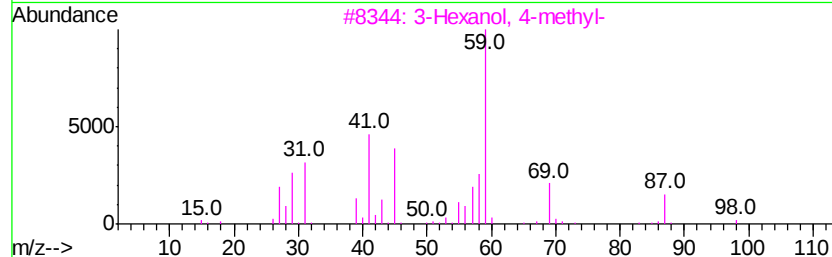
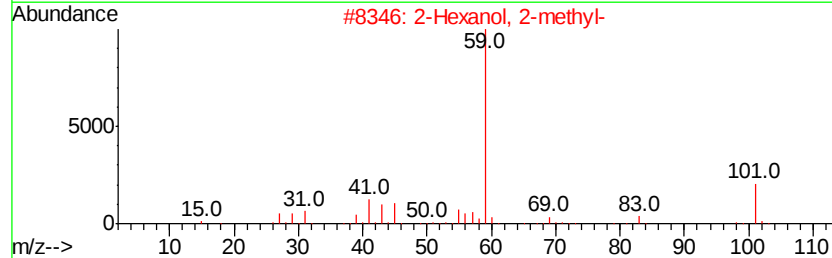
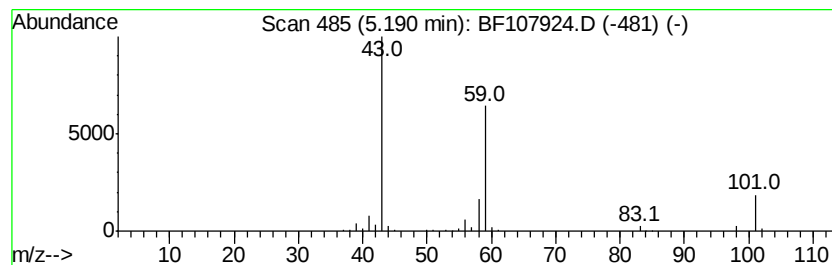
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.19	9.93 ng	393680	1,4-Dichlorobenzene-d4	6.96

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2		2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	38
3		3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33
4		1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10
5		Acetone	58	C3H6O	000067-64-1	9



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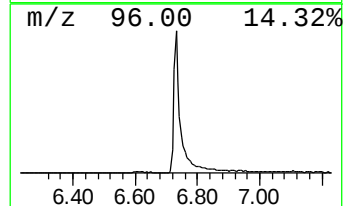
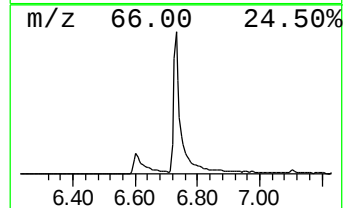
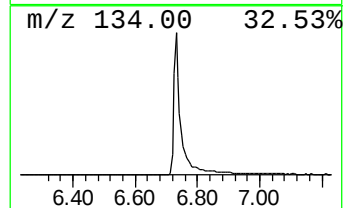
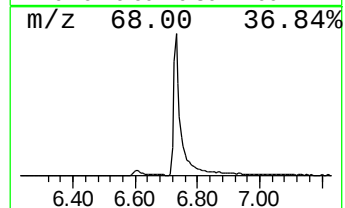
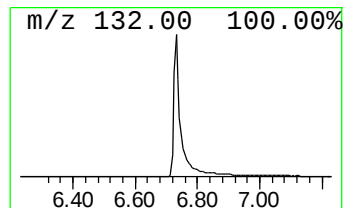
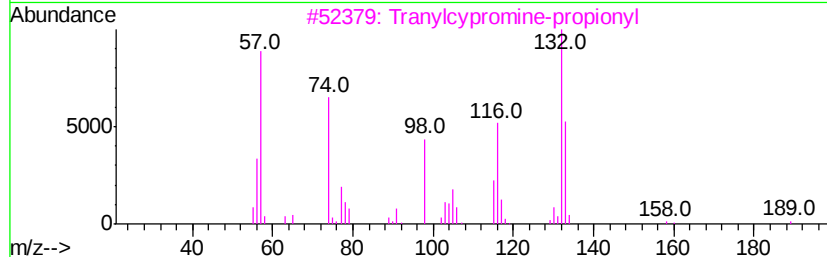
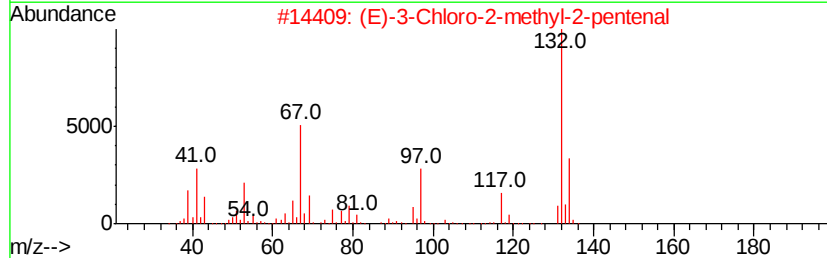
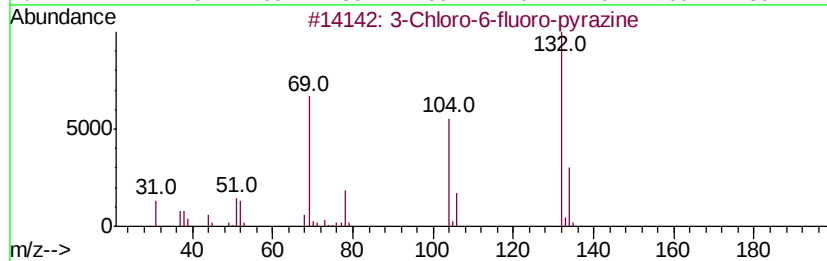
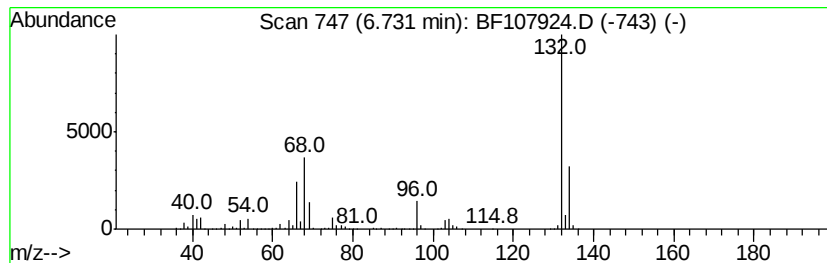
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 2 unknown6.73 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.73	77.20 ng	3061580	1,4-Dichlorobenzene-d4	6.96

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	35
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	25
3		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	12
4		4-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-60-2	9
5		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9



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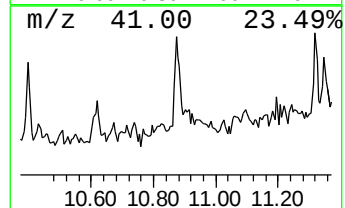
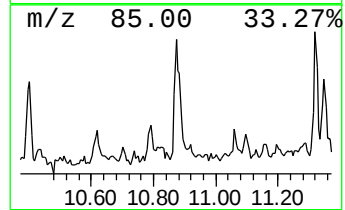
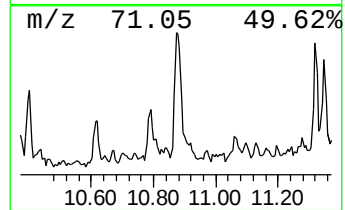
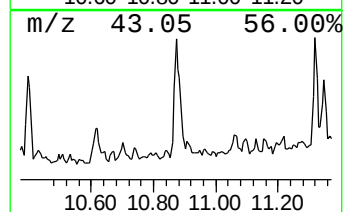
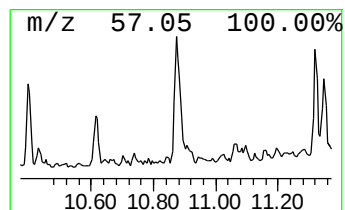
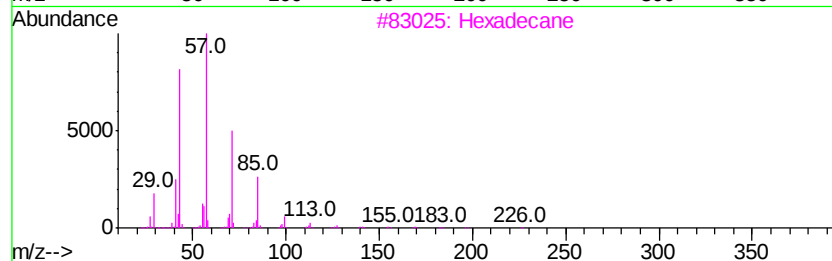
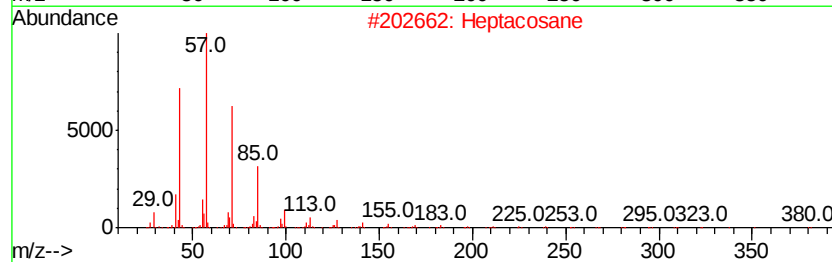
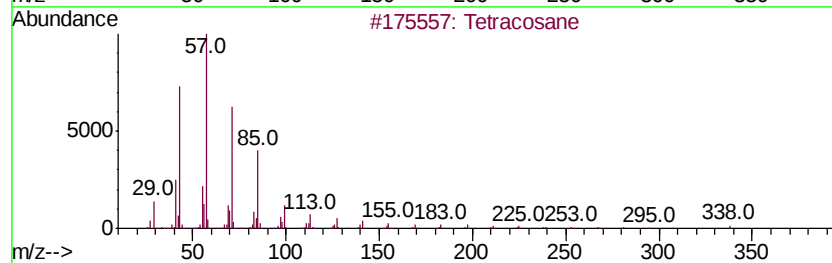
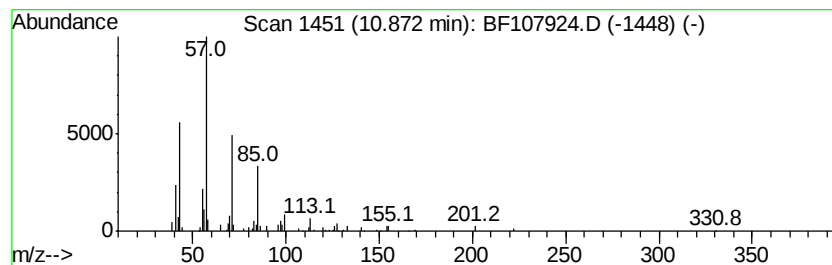
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 4 Tetracosane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.87	2.58 ng	108512	Phenanthrene-d10	11.49

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetracosane	338	C24H50	000646-31-1	80
2			Heptacosane	380	C27H56	000593-49-7	80
3			Hexadecane	226	C16H34	000544-76-3	78
4			Tridecane, 1-iodo-	310	C13H27I	035599-77-0	72
5			triacontane	422	C30H62	000638-68-6	72



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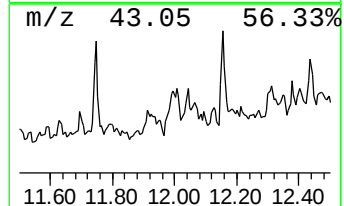
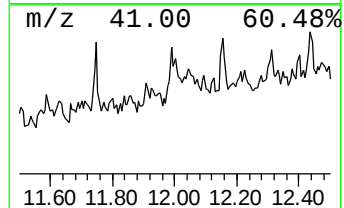
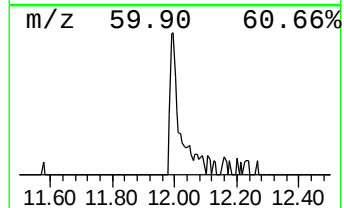
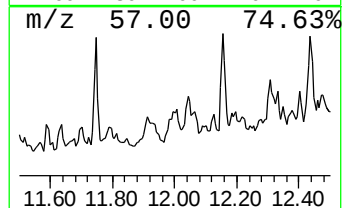
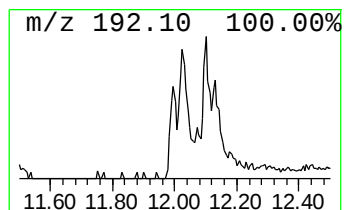
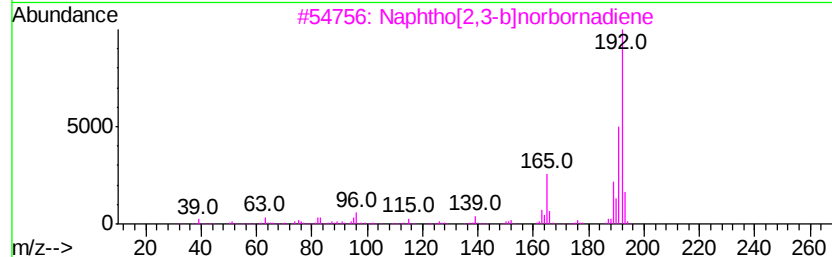
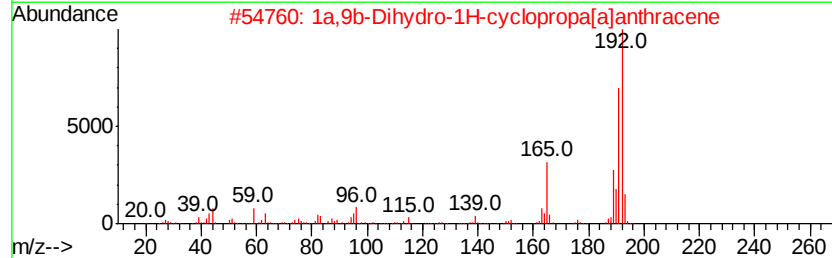
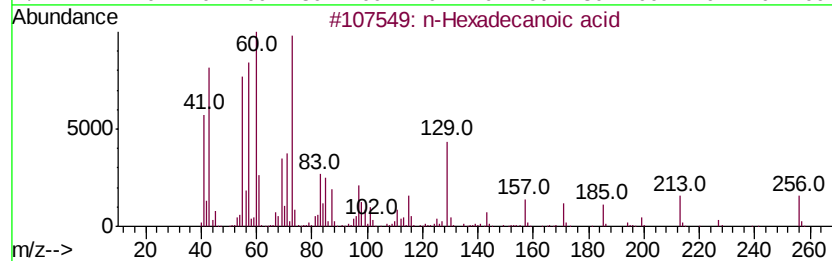
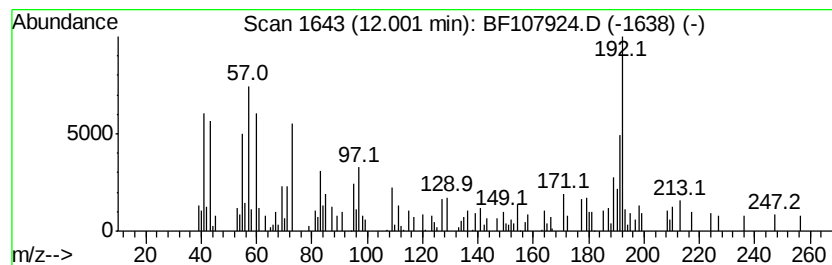
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 n-Hexadecanoic acid Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.00	2.84 ng	119710	Phenanthrene-d10	11.49

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	55
2		1a,9b-Dihydro-1H-cyclopropa[an...	192	C15H12	006109-24-6	38
3		Naphtho[2,3-b]norbornadiene	192	C15H12	107426-38-0	38
4		Anthracene, 1-methyl-	192	C15H12	000610-48-0	35
5		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	35



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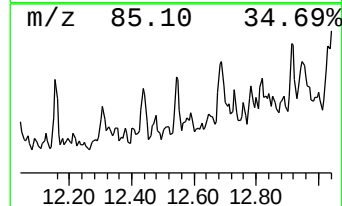
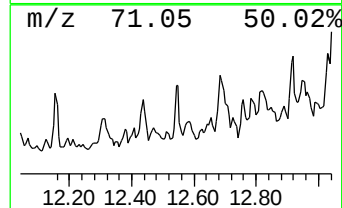
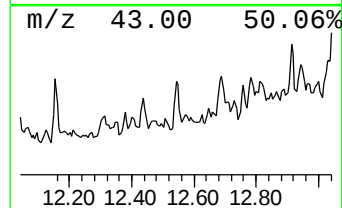
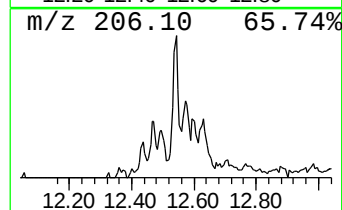
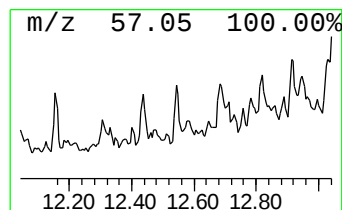
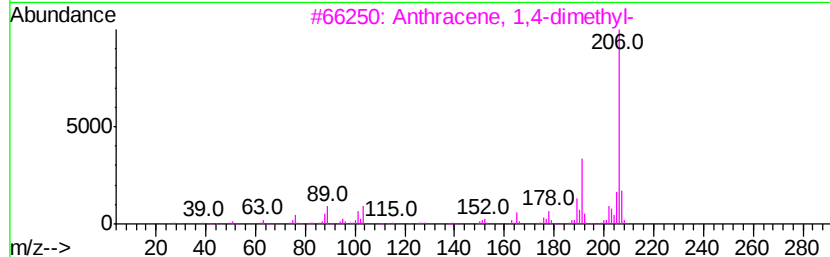
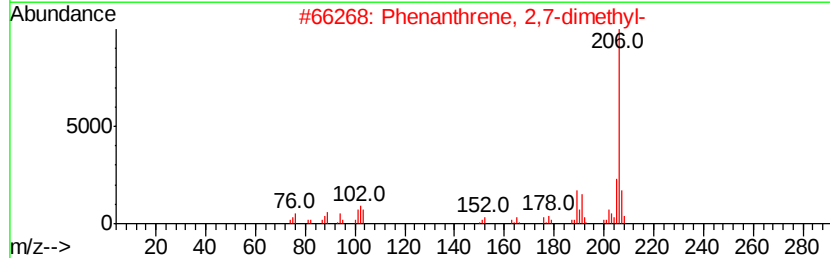
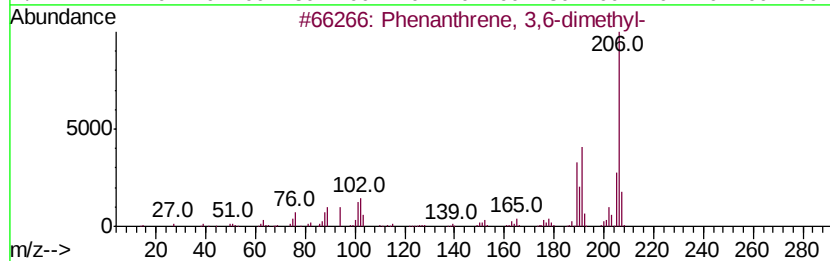
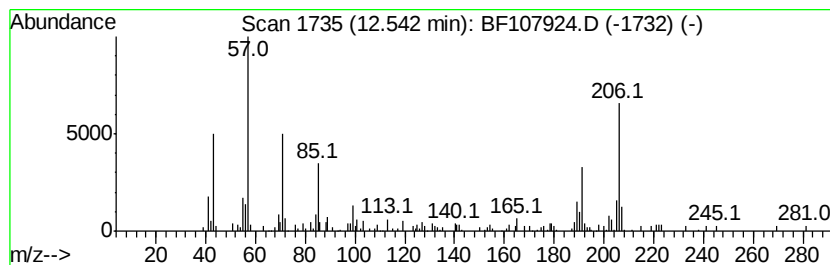
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Phenanthrene, 3,6-dimethyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.54	2.18 ng	92010	Phenanthrene-d10	11.49	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 3,6-dimethyl-	206	C16H14	001576-67-6	87
2	Phenanthrene, 2,7-dimethyl-	206	C16H14	001576-69-8	86
3	Anthracene, 1,4-dimethyl-	206	C16H14	000781-92-0	78
4	Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	70
5	Naphthalene, 1,2-dihydro-1-phenyl-	206	C16H14	016606-46-5	58



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF080218\
Data File : BF107924.D
Acq On : 2 Aug 2018 22:01
Operator : JU/SJ
Sample : J4285-21
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
KG-PIT-6

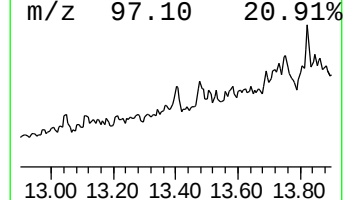
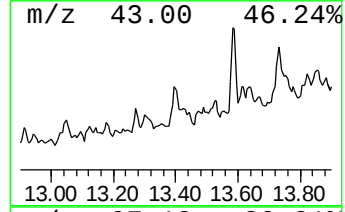
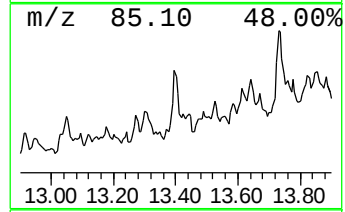
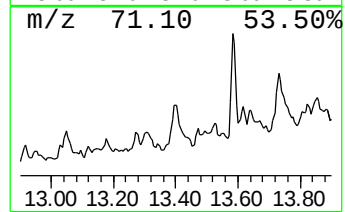
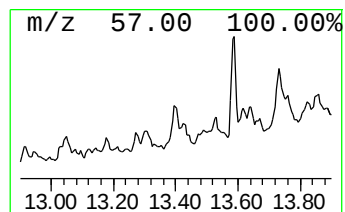
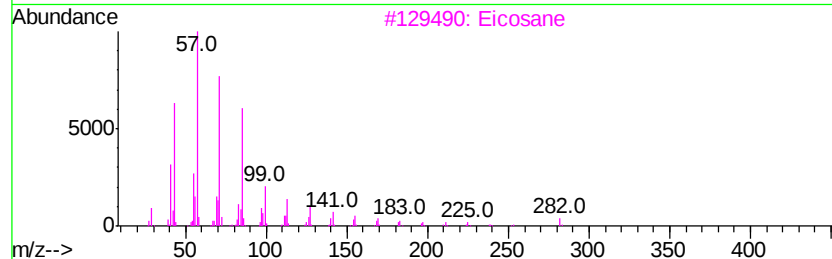
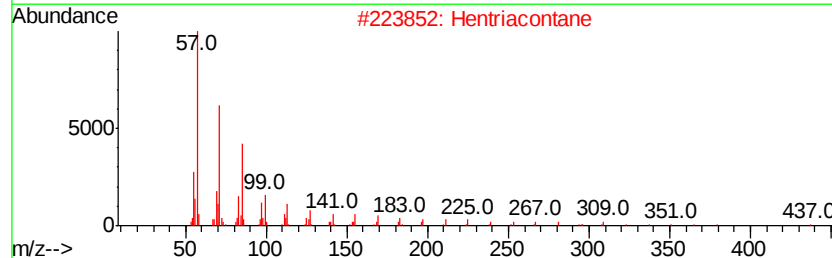
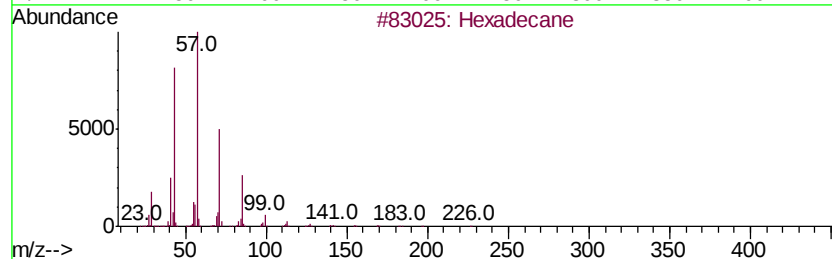
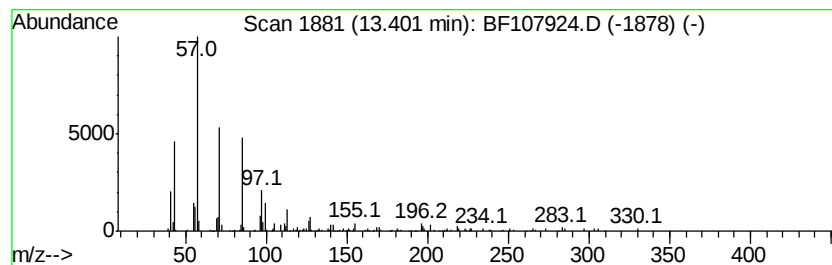
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Hexadecane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.40	2.49 ng	129177	Chrysene-d12	14.14

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadecane	226	C16H34	000544-76-3	91
2		Hentriacontane	437	C31H64	000630-04-6	78
3		Eicosane	282	C20H42	000112-95-8	72
4		Heneicosane	296	C21H44	000629-94-7	72
5		Decane, 3,8-dimethyl-	170	C12H26	017312-55-9	72



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF080218\
Data File : BF107924.D
Acq On : 2 Aug 2018 22:01
Operator : JU/SJ
Sample : J4285-21
Misc :
ALS Vial : 20 Sample Multiplier: 1

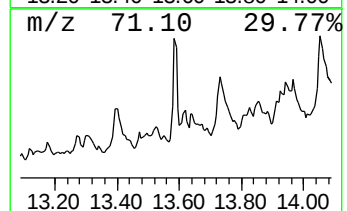
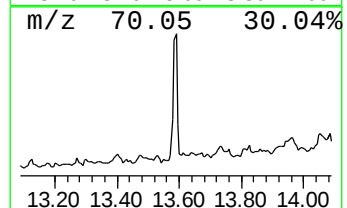
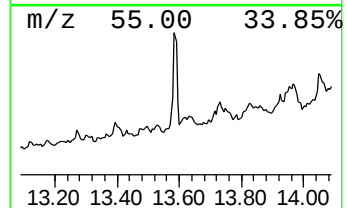
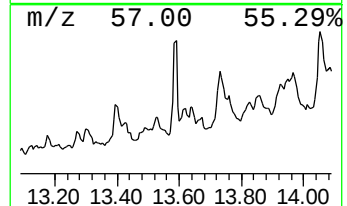
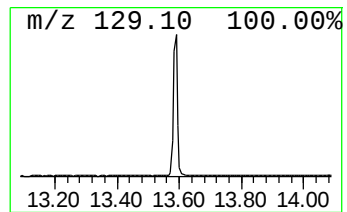
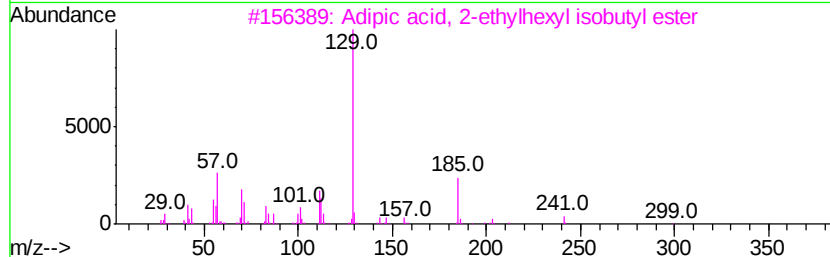
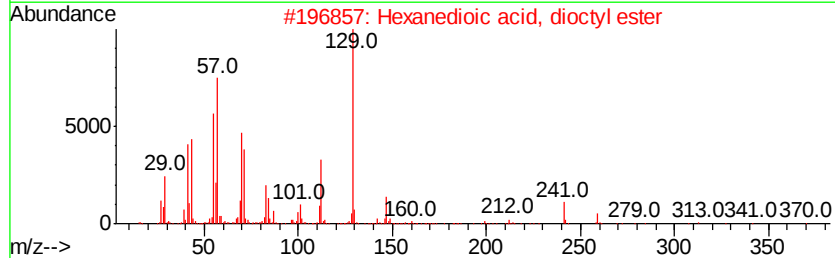
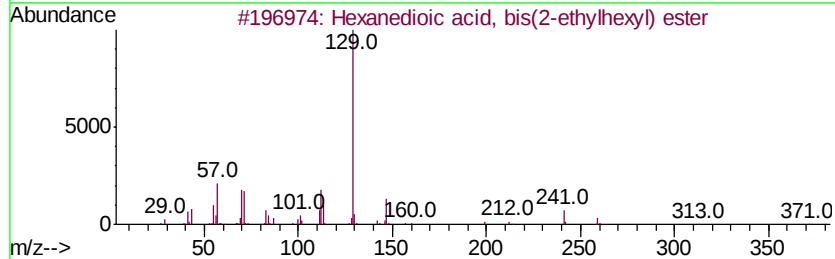
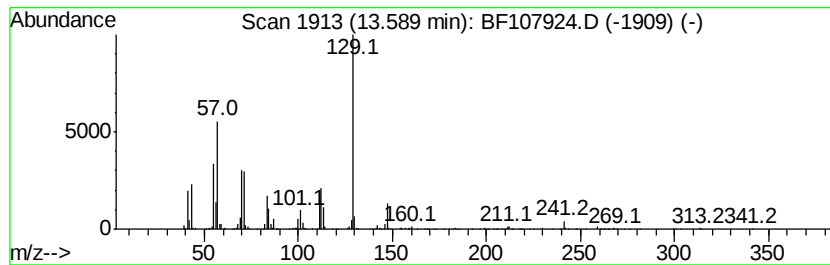
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Hexanedioic acid, bis(2-eth... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.		
13.59	13.46 ng	696836	Chrysene-d12	14.14		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Hexanedioic acid, bis(2-ethylhex...	370	C22H42O4	000103-23-1	91
2		Hexanedioic acid, dioctyl ester	370	C22H42O4	000123-79-5	70
3		Adipic acid, 2-ethylhexyl isobut...	314	C18H34O4	1000324-48-0	64
4		Adipic acid, 2-ethylhexyl isohex...	342	C20H38O4	1000324-48-3	53
5		Adipic acid, cycloheptyl isohexy...	326	C19H34O4	1000324-71-8	50



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF080218\
Data File : BF107924.D
Acq On : 2 Aug 2018 22:01
Operator : JU/SJ
Sample : J4285-21
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
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ClientSampleId :
KG-PIT-6

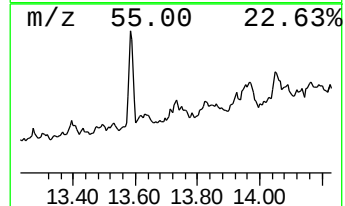
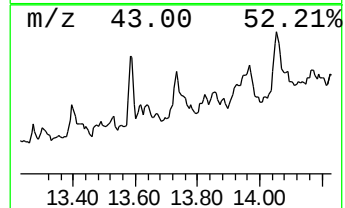
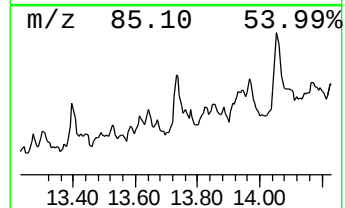
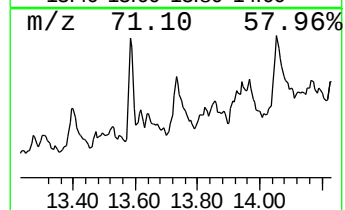
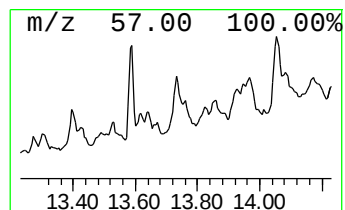
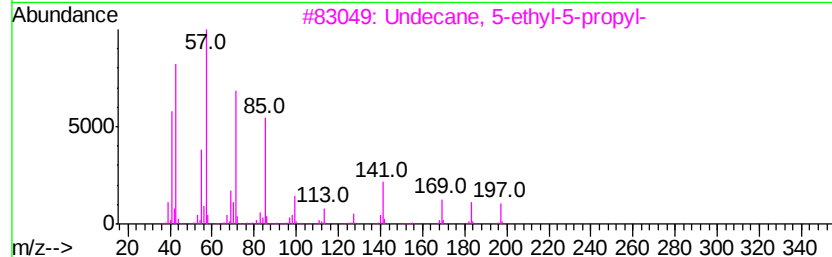
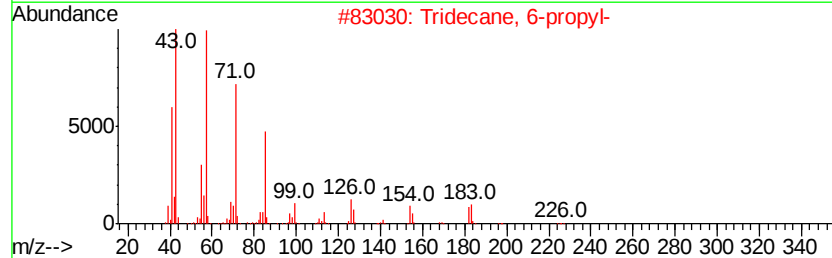
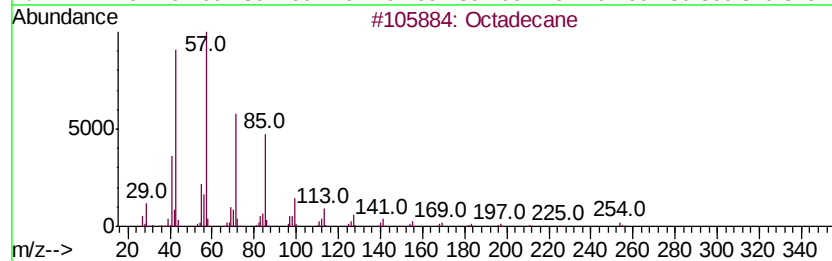
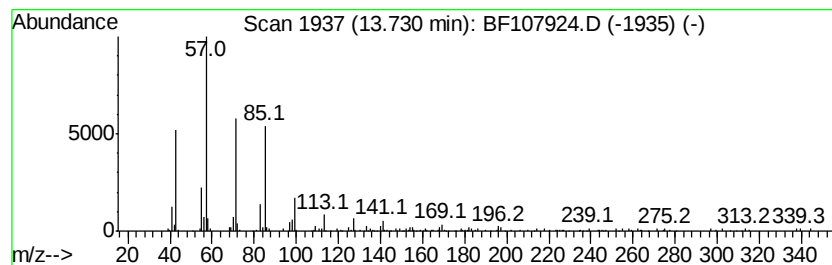
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Octadecane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.73	4.37 ng	226376	Chrysene-d12	14.14

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecane	254	C18H38	000593-45-3	80
2		Tridecane, 6-propyl-	226	C16H34	055045-10-8	80
3		Undecane, 5-ethyl-5-propyl-	226	C16H34	002755-07-9	72
4		Heptane, 2,2,3,3,5,6,6-heptamethyl-	198	C14H30	007225-67-4	53
5		Sulfurous acid, butyl octadecyl ...	390	C22H46O3S	1000309-18-5	53



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF080218\
Data File : BF107924.D
Acq On : 2 Aug 2018 22:01
Operator : JU/SJ
Sample : J4285-21
Misc :
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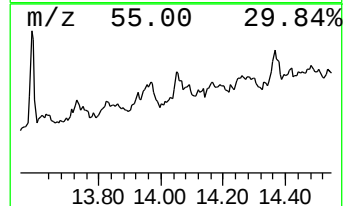
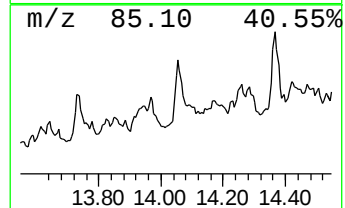
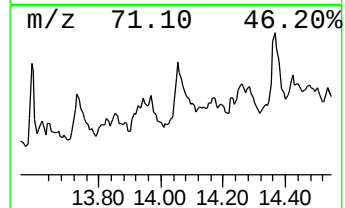
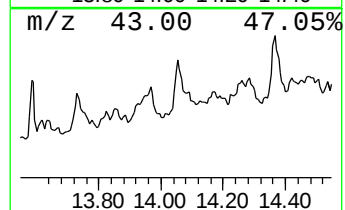
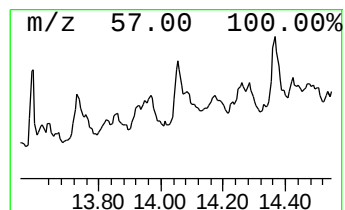
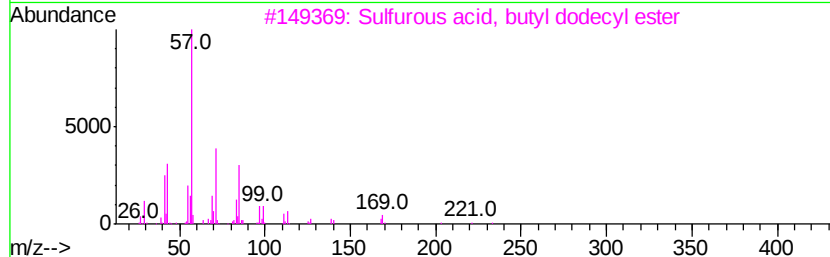
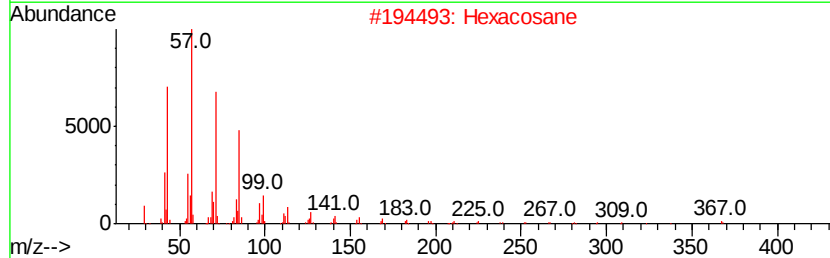
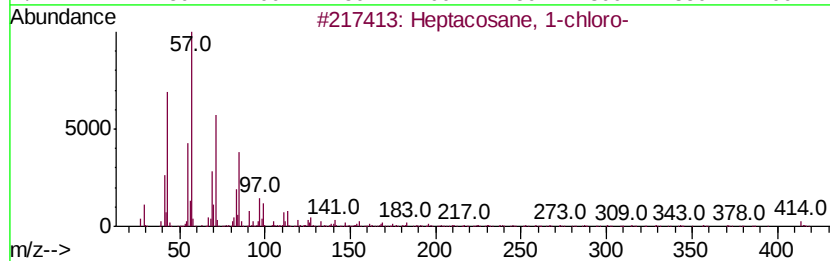
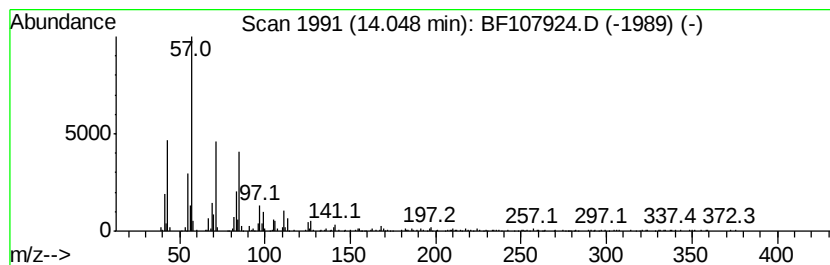
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Heptacosane, 1-chloro- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.05	9.28 ng	480337	Chrysene-d12	14.14

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptacosane, 1-chloro-	414	C27H55Cl	062016-79-9	87
2		Hexacosane	366	C26H54	000630-01-3	87
3		Sulfurous acid, butyl dodecyl ester	306	C16H34O3S	1000309-17-9	86
4		Pentacosane, 13-undecyl-	507	C36H74	055517-89-0	86
5		2-methyloctacosane	408	C29H60	1000376-72-8	86



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF080218\
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Acq On : 2 Aug 2018 22:01
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Sample : J4285-21
Misc :
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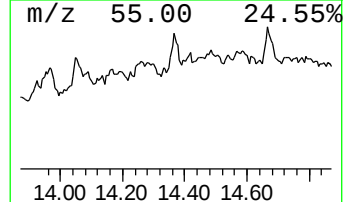
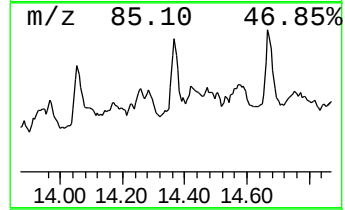
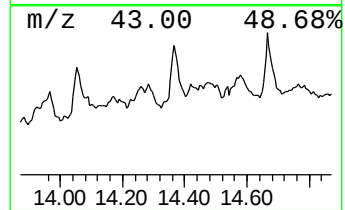
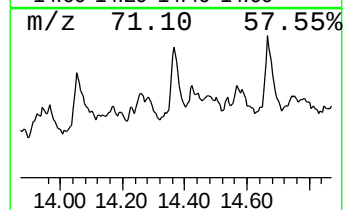
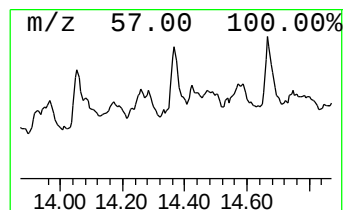
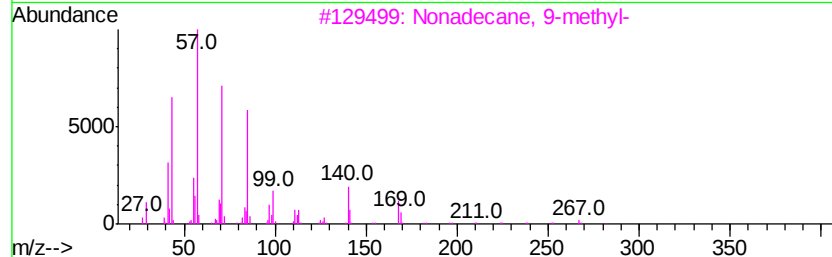
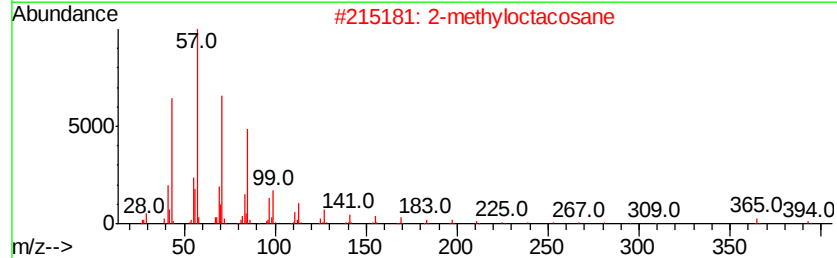
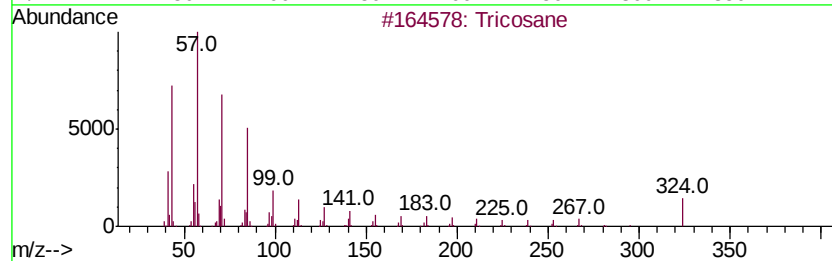
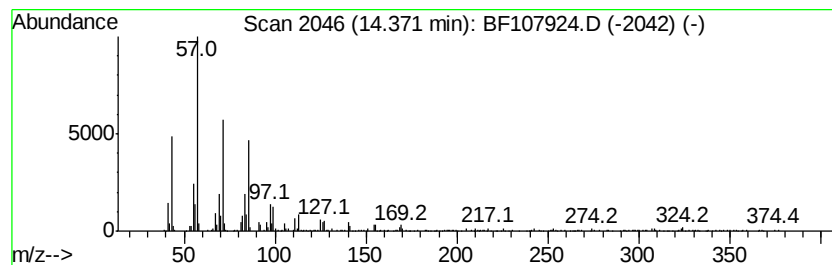
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF073018.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Tricosane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD		R.T.
14.37	10.87 ng	563161	Chrysene-d12		14.14
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tricosane	324	C23H48	000638-67-5	91
2	2-methyloctacosane	408	C29H60	1000376-72-8	86
3	Nonadecane, 9-methyl-	282	C20H42	013287-24-6	86
4	Hentriacontane	437	C31H64	000630-04-6	86
5	Docosane, 11-butyl-	366	C26H54	013475-76-8	86



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF080218\
Data File : BF107924.D
Acq On : 2 Aug 2018 22:01
Operator : JU/SJ
Sample : J4285-21
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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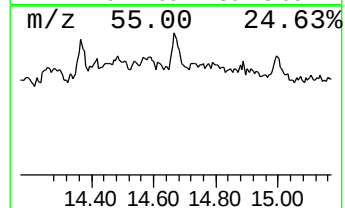
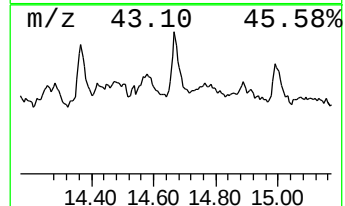
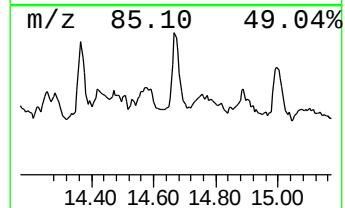
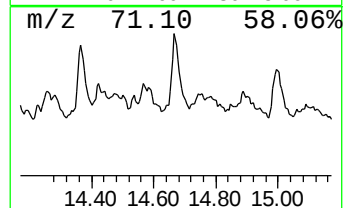
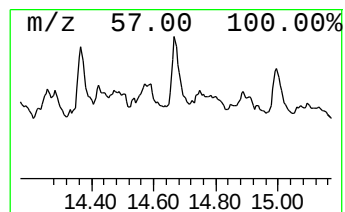
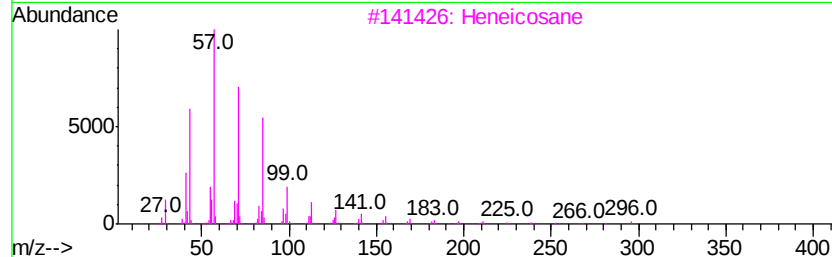
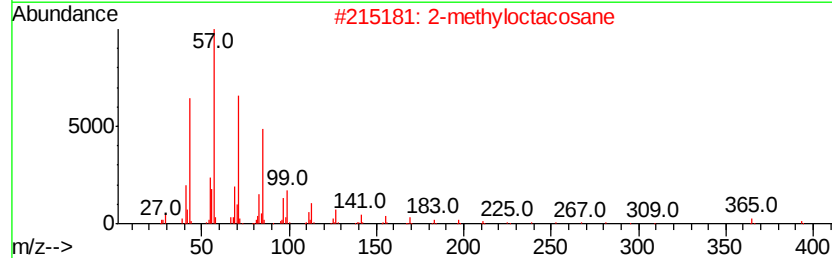
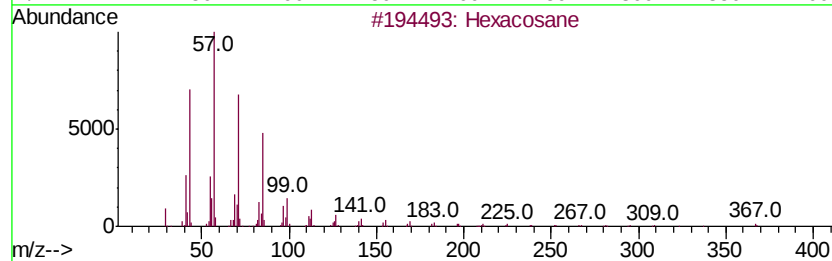
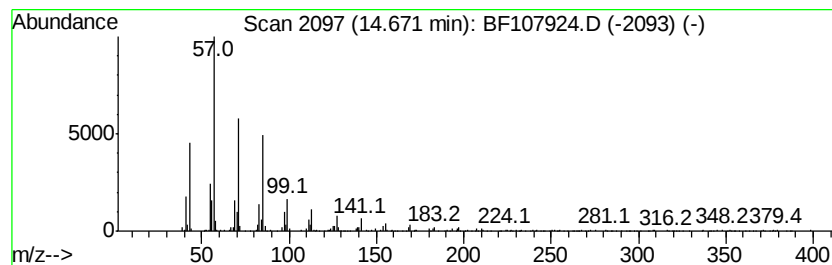
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Hexacosane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.67	10.96 ng	567656	Chrysene-d12	14.14

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexacosane	366	C26H54	000630-01-3	91
2		2-methyloctacosane	408	C29H60	1000376-72-8	90
3		Heneicosane	296	C21H44	000629-94-7	90
4		Triacontane	422	C30H62	000638-68-6	87
5		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF080218\
Data File : BF107924.D
Acq On : 2 Aug 2018 22:01
Operator : JU/SJ
Sample : J4285-21
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
KG-PIT-6

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF073018.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT 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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unknown6.73	6.73	77.2	ng	3061580	1	6.96	793134	20.0
Tetracosane	10.87	2.6	ng	108512	4	11.49	842593	20.0
n-Hexadecanoic acid	12.00	2.8	ng	119710	4	11.49	842593	20.0
Phenanthrene, 3,6...	12.54	2.2	ng	92010	4	11.49	842593	20.0
Hexadecane	13.40	2.5	ng	129177	5	14.14	1035730	20.0
Hexanedioic acid,...	13.59	13.5	ng	696836	5	14.14	1035730	20.0
Octadecane	13.73	4.4	ng	226376	5	14.14	1035730	20.0
Heptacosane, 1-ch...	14.05	9.3	ng	480337	5	14.14	1035730	20.0
Tricosane	14.37	10.9	ng	563161	5	14.14	1035730	20.0
Hexacosane	14.67	11.0	ng	567656	5	14.14	1035730	20.0