

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF101018\
 Data File : BF109747.D
 Acq On : 10 Oct 2018 17:27
 Operator : JU/SJ
 Sample : J5369-01
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CONCRETE-PILE

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-BF100818.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	4.240	360	366	386	rBV	859311	1305057	58.16%	8.023%
2	4.893	470	477	480	rBV	1019783	1367745	60.96%	8.409%
3	5.316	546	549	563	rBV	78974	158331	7.06%	0.973%
4	6.340	720	723	741	rBV	683552	1024911	45.68%	6.301%
5	6.463	741	744	768	rVB	391358	561685	25.03%	3.453%
6	6.693	780	783	806	rBV	264122	422924	18.85%	2.600%
7	6.851	806	810	837	rVV	1112315	1262847	56.28%	7.764%
8	7.022	837	839	856	rVB4	11309	27848	1.24%	0.171%
9	7.257	876	879	908	rBV	696677	1023646	45.62%	6.293%
10	7.975	997	1001	1024	rBV	418629	577024	25.72%	3.547%
11	9.051	1180	1184	1197	rBV	1686107	1997203	89.01%	12.279%
12	9.139	1197	1199	1210	rVB3	24788	48402	2.16%	0.298%
13	9.445	1247	1251	1269	rBV	184628	231554	10.32%	1.424%
14	9.722	1294	1298	1316	rBV	652181	677054	30.17%	4.162%
15	11.198	1546	1549	1568	rBV	545295	773630	34.48%	4.756%
16	12.663	1795	1798	1800	rBV	29578	24458	1.09%	0.150%
17	12.686	1800	1802	1806	rVB	34094	28631	1.28%	0.176%
18	12.780	1814	1818	1831	rBV	2203277	2243770	100.00%	13.794%
19	13.016	1853	1858	1869	rVB	51128	54002	2.41%	0.332%
20	13.357	1912	1916	1919	rBV	78997	76715	3.42%	0.472%
21	13.686	1967	1972	1982	rBV	150467	222474	9.92%	1.368%
22	13.827	1992	1996	2005	rBV	613658	708067	31.56%	4.353%
23	13.998	2022	2025	2028	rBV	115741	101867	4.54%	0.626%
24	14.304	2073	2077	2081	rBV	105947	105218	4.69%	0.647%
25	14.592	2123	2126	2131	rBV	113925	117109	5.22%	0.720%
26	14.663	2134	2138	2144	rVV	166026	163157	7.27%	1.003%
27	14.904	2175	2179	2184	rBV	97328	109868	4.90%	0.675%
28	15.227	2229	2234	2250	rVB	340862	645246	28.76%	3.967%
29	15.645	2300	2305	2310	rBV	64315	92805	4.14%	0.571%
30	16.104	2378	2383	2389	rVB4	45556	77729	3.46%	0.478%
31	16.633	2468	2473	2481	rVB3	18313	34739	1.55%	0.214%

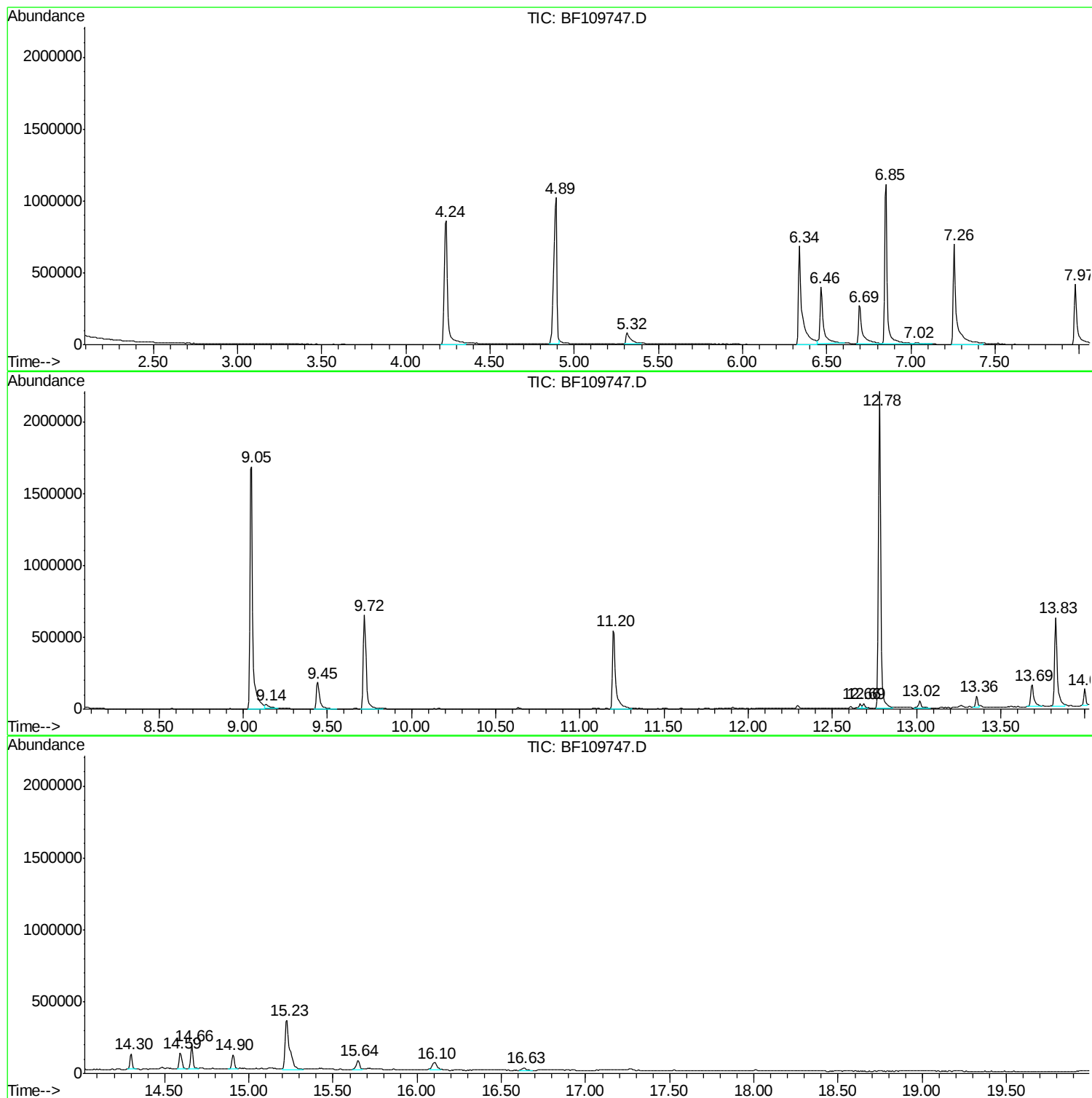
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF100818.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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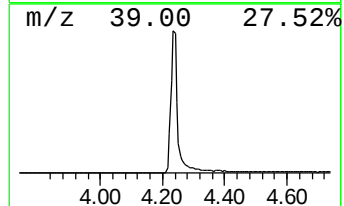
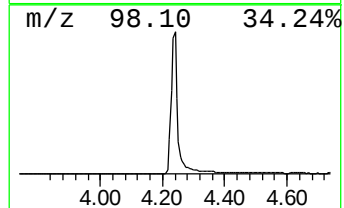
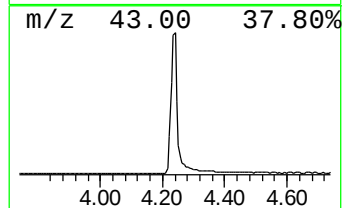
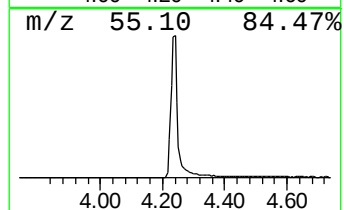
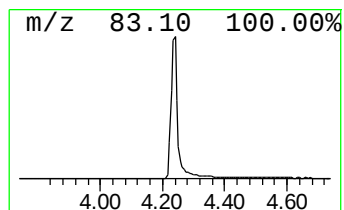
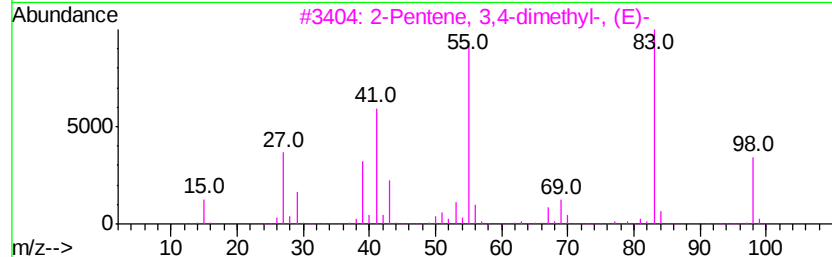
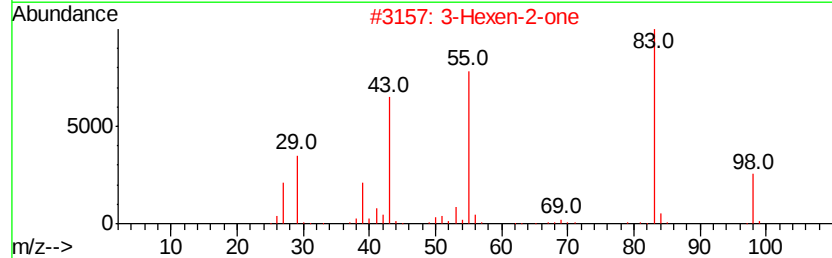
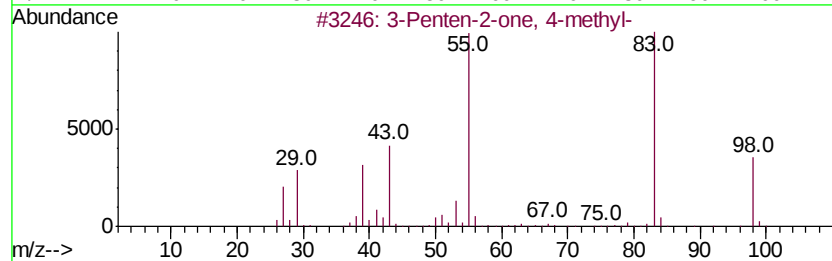
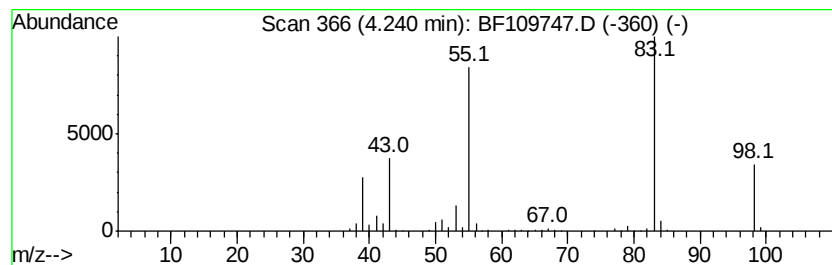
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 3-Penten-2-one, 4-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.24	61.72 ng	1305060	1,4-Dichlorobenzene-d4	6.70

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Penten-2-one, 4-methyl-	98	C6H10O	000141-79-7	95
2		3-Hexen-2-one	98	C6H10O	000763-93-9	91
3		2-Pentene, 3,4-dimethyl-, (E)-	98	C7H14	004914-92-5	90
4		2-Pentene, 3,4-dimethyl-, (Z)-	98	C7H14	004914-91-4	86
5		2-Pentene, 2,4-dimethyl-	98	C7H14	000625-65-0	80



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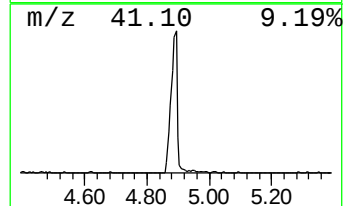
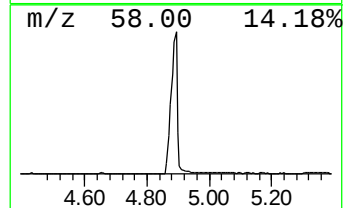
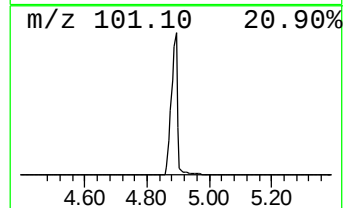
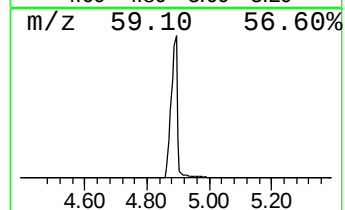
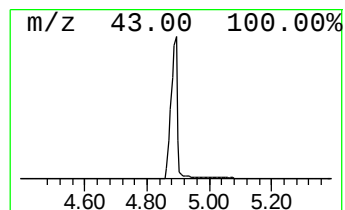
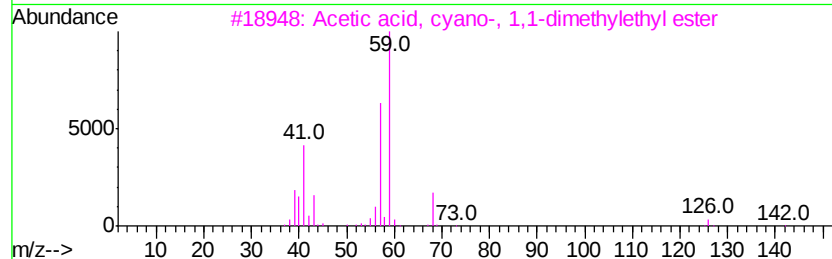
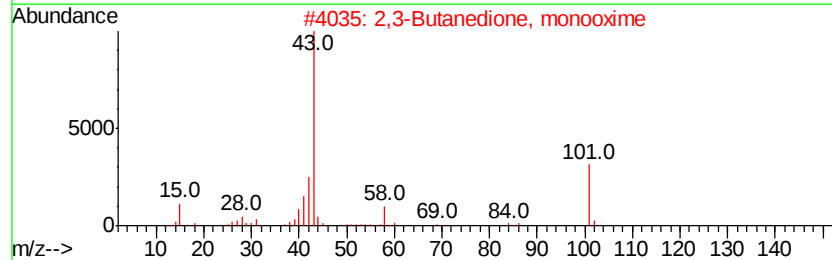
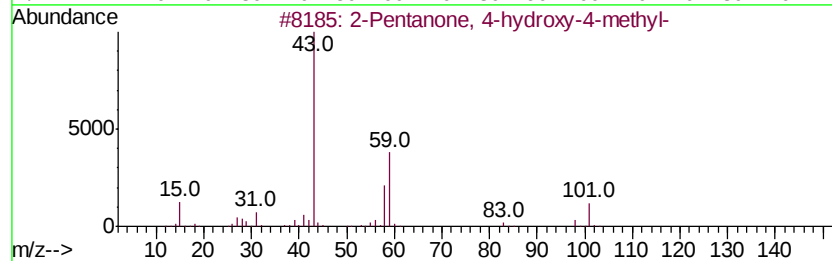
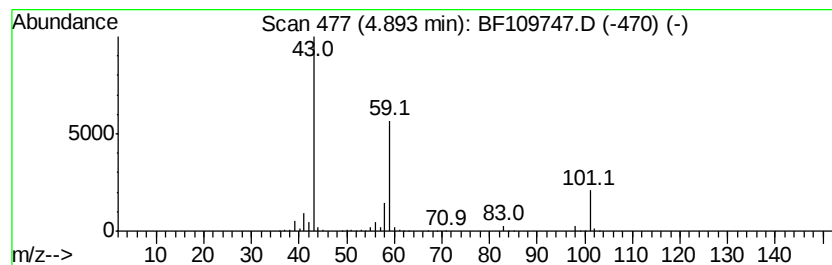
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.89	64.68 ng	1367750	1,4-Dichlorobenzene-d4	6.70

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	17
3		Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	17
4		5-Hexen-2-one	98	C6H10O	000109-49-9	9
5		Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9



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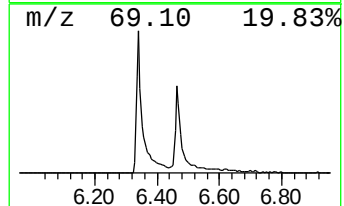
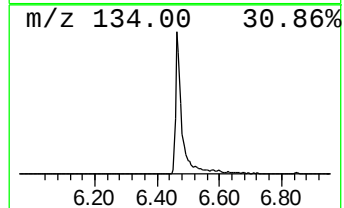
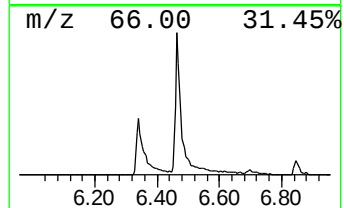
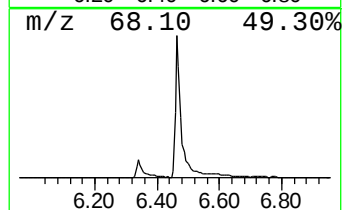
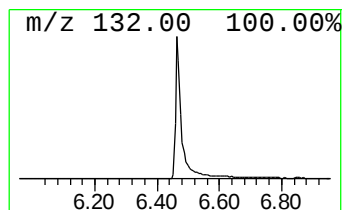
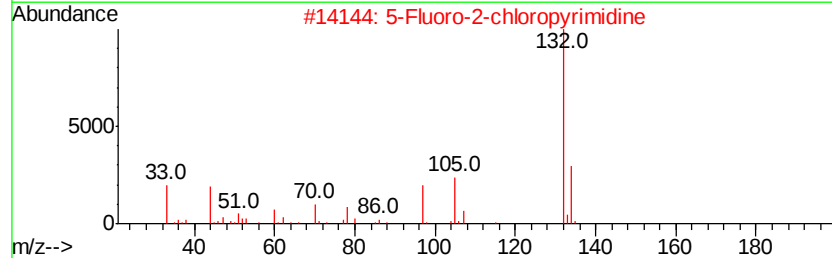
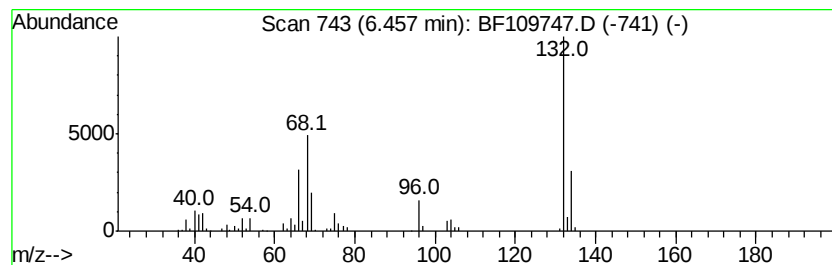
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 unknown6.46 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.46	26.56 ng	561685	1,4-Dichlorobenzene-d4	6.70

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27	
2	5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12	
3	Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	10	
4	4-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-60-2	9	
5	1-Aminoindan	133	C9H11N	034698-41-4	9	



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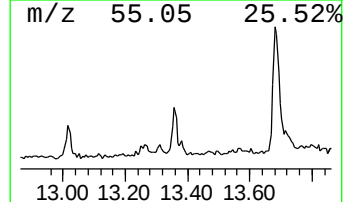
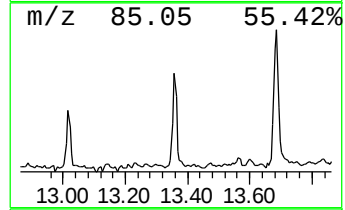
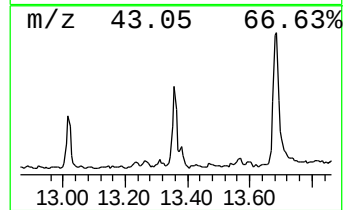
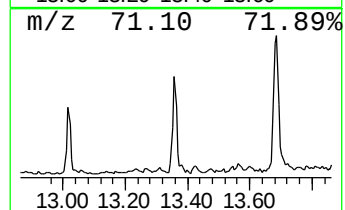
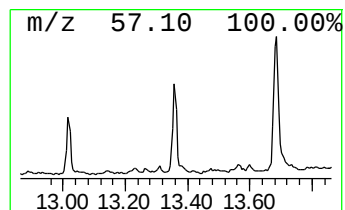
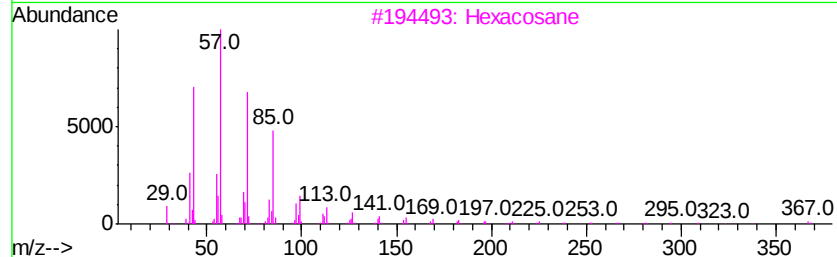
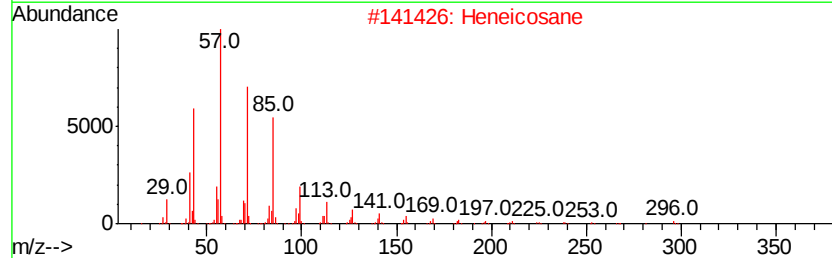
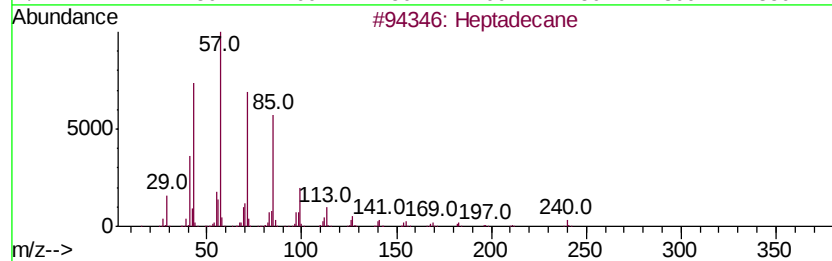
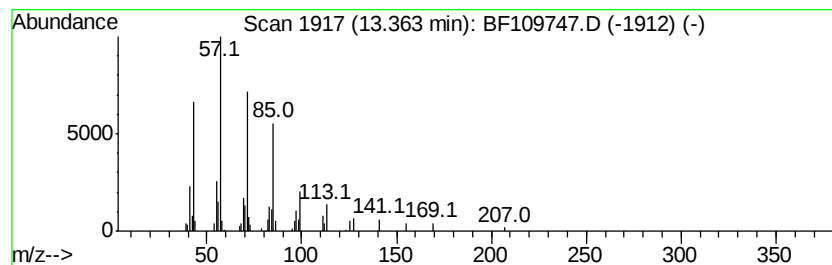
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Peak Number 5 Heptadecane Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.36	2.17 ng	76715	Chrysene-d12	13.83

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane		240	C17H36	000629-78-7	91
2	Heneicosane		296	C21H44	000629-94-7	91
3	Hexacosane		366	C26H54	000630-01-3	90
4	Octadecane		254	C18H38	000593-45-3	90
5	Hexadecane		226	C16H34	000544-76-3	86



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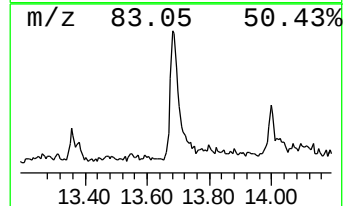
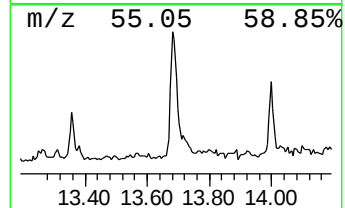
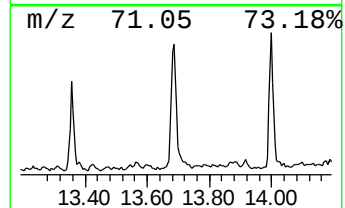
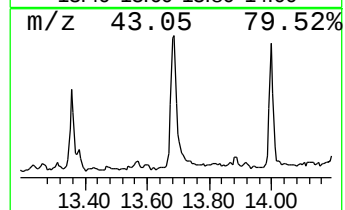
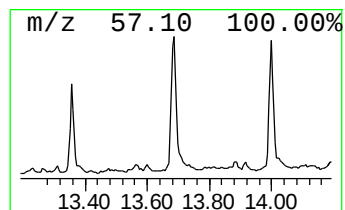
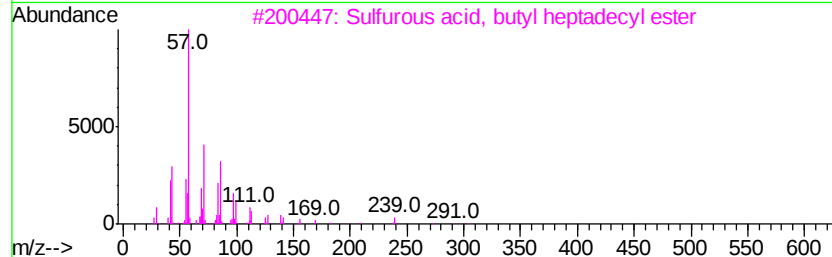
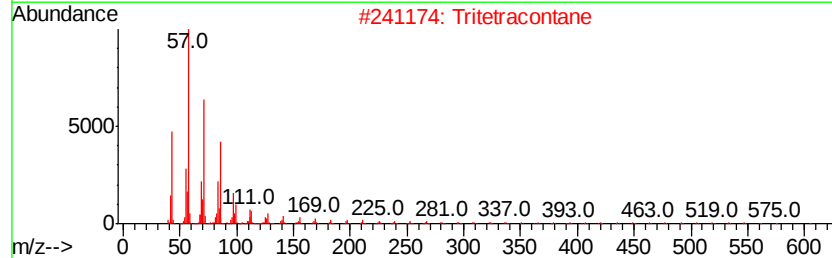
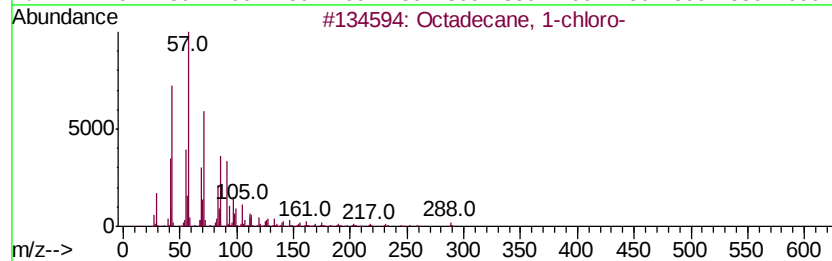
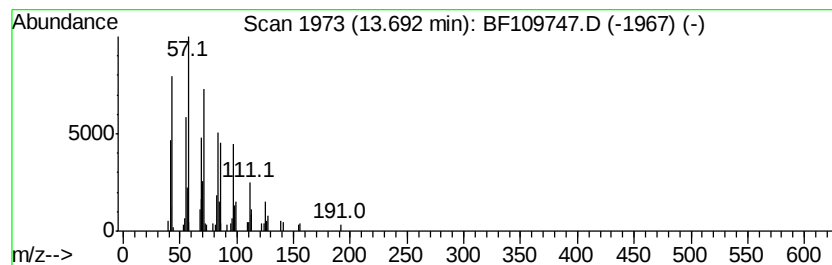
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 6 Octadecane, 1-chloro- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.69	6.28 ng	222474	Chrysene-d12	13.83

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecane, 1-chloro-	288	C18H37Cl	003386-33-2	90
2		Tritetracontane	605	C43H88	007098-21-7	86
3		Sulfurous acid, butyl heptadecyl...	376	C21H44O3S	1000309-18-4	72
4		Hexacosane	366	C26H54	000630-01-3	72
5		Hexadecane	226	C16H34	000544-76-3	64



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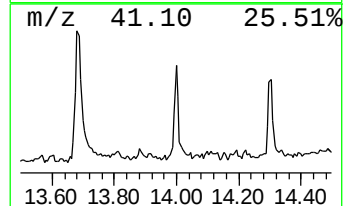
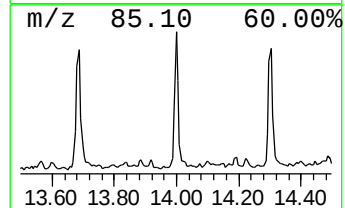
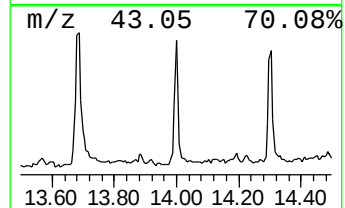
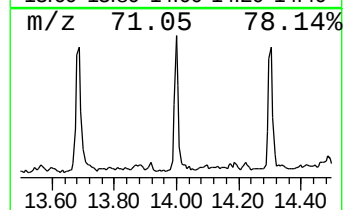
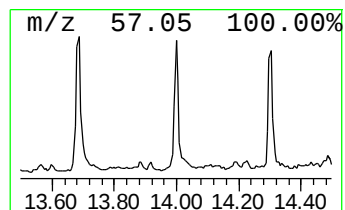
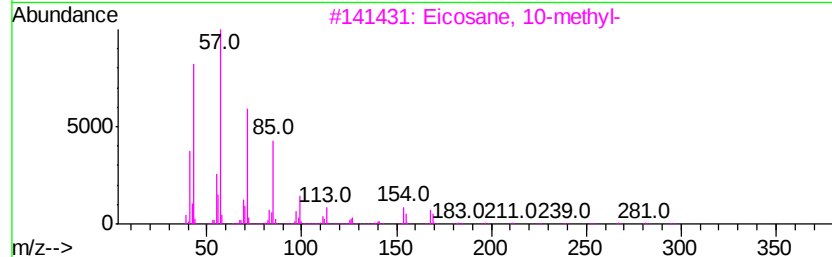
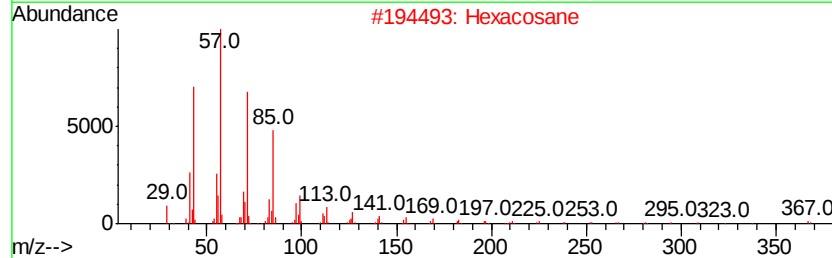
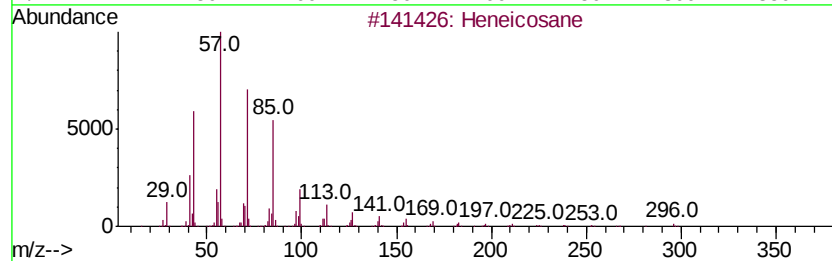
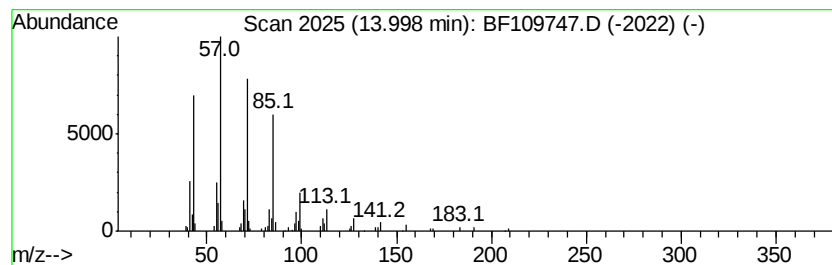
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Heneicosane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.00	2.88 ng	101867	Chrysene-d12	13.83

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane		296	C21H44	000629-94-7	91
2	Hexacosane		366	C26H54	000630-01-3	90
3	Eicosane, 10-methyl-		296	C21H44	054833-23-7	86
4	Hexadecane		226	C16H34	000544-76-3	86
5	Tetracosane		338	C24H50	000646-31-1	86



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF101018\
Data File : BF109747.D
Acq On : 10 Oct 2018 17:27
Operator : JU/SJ
Sample : J5369-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
CONCRETE-PILE

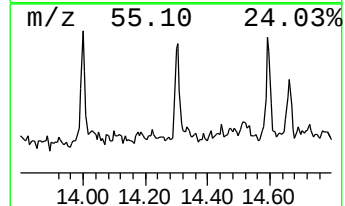
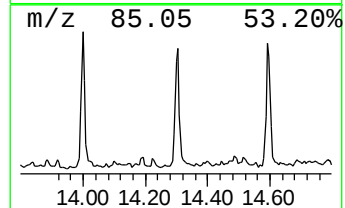
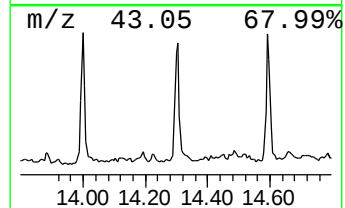
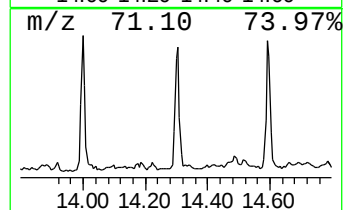
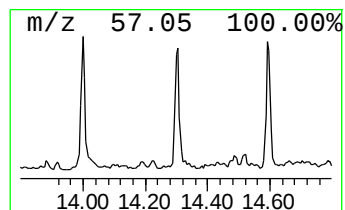
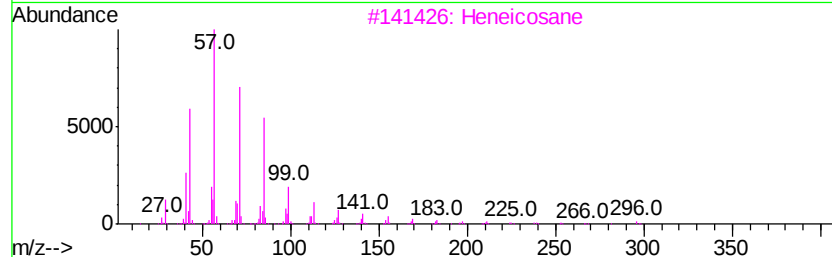
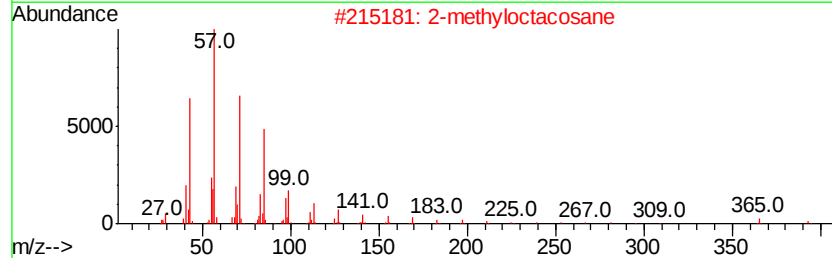
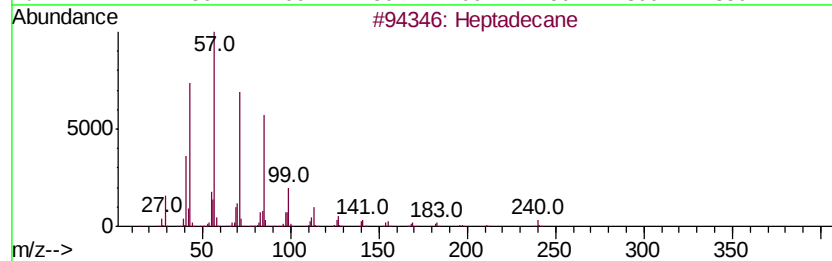
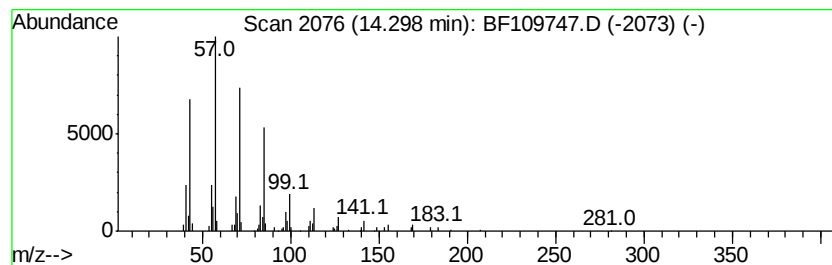
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 2-methyloctacosane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.30	2.97 ng	105218	Chrysene-d12	13.83

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heptadecane	240	C17H36	000629-78-7	91
2		2-methyloctacosane	408	C29H60	1000376-72-8	91
3		Heneicosane	296	C21H44	000629-94-7	91
4		Hexacosane	366	C26H54	000630-01-3	90
5		Tetracosane	338	C24H50	000646-31-1	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF101018\
Data File : BF109747.D
Acq On : 10 Oct 2018 17:27
Operator : JU/SJ
Sample : J5369-01
Misc :
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Instrument :
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ClientSampleId :
CONCRETE-PILE

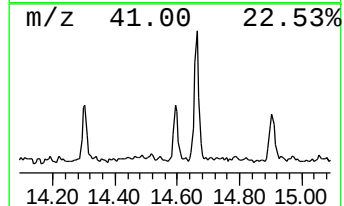
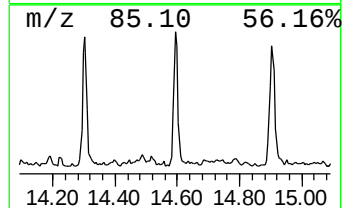
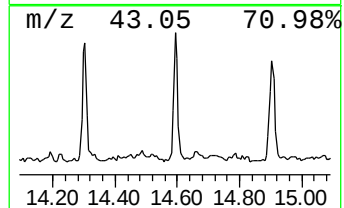
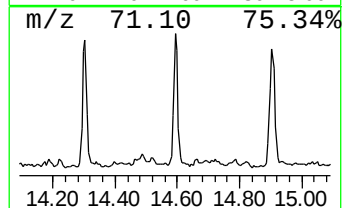
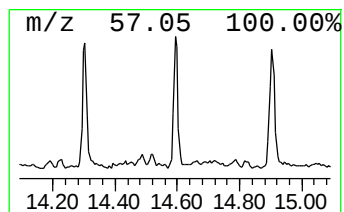
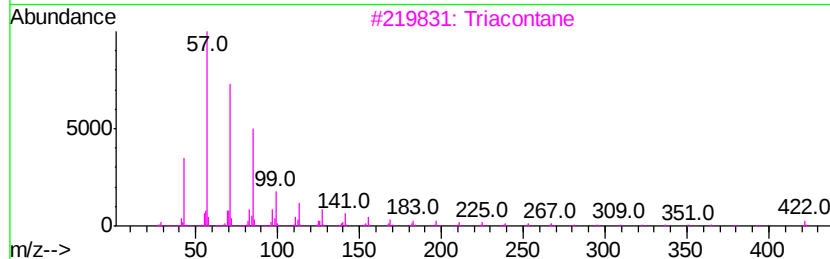
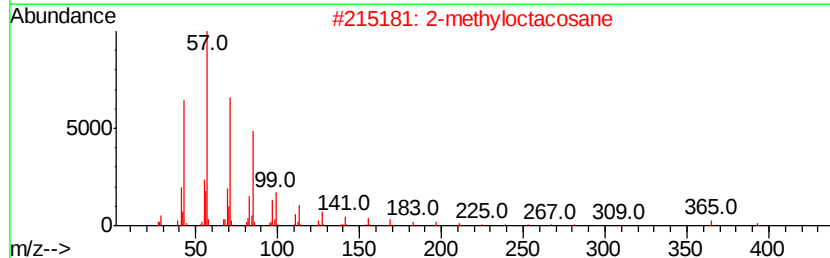
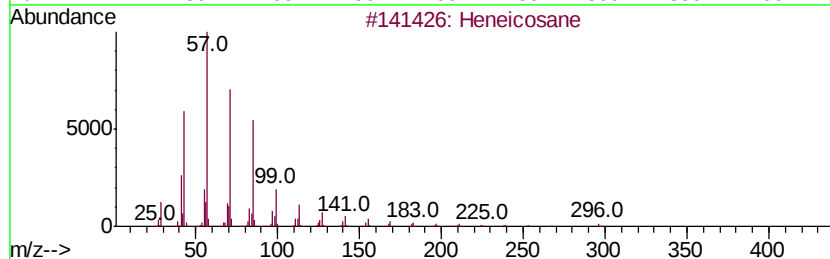
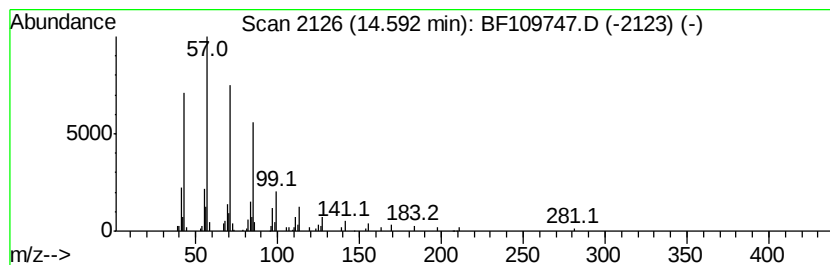
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Triacontane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.59	3.63 ng	117109	Perylene-d12	15.23

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



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF101018\
Data File : BF109747.D
Acq On : 10 Oct 2018 17:27
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Misc :
ALS Vial : 11 Sample Multiplier: 1

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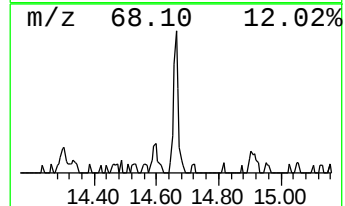
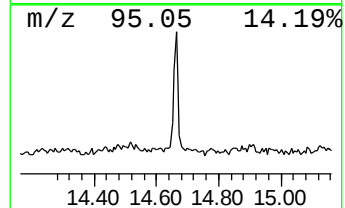
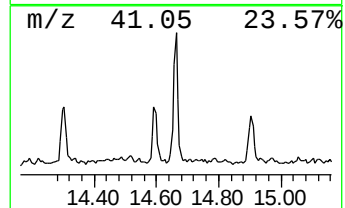
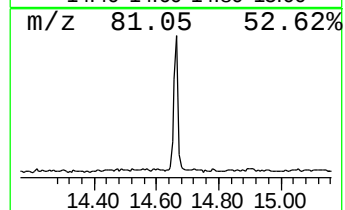
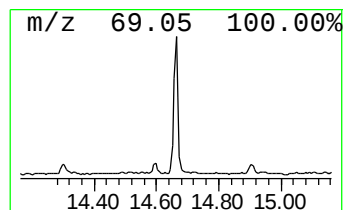
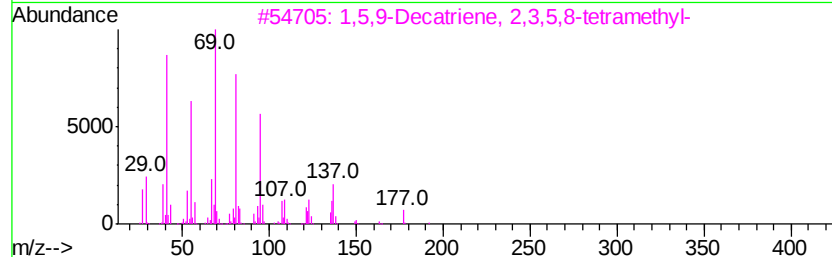
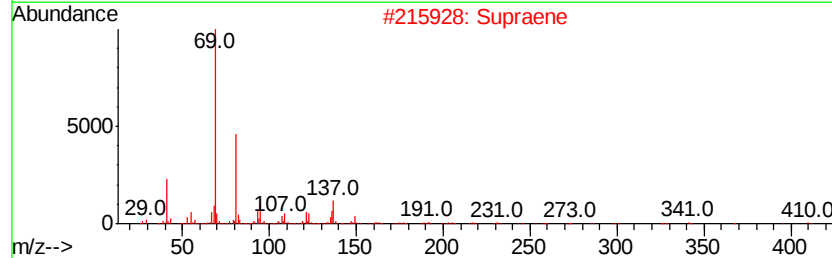
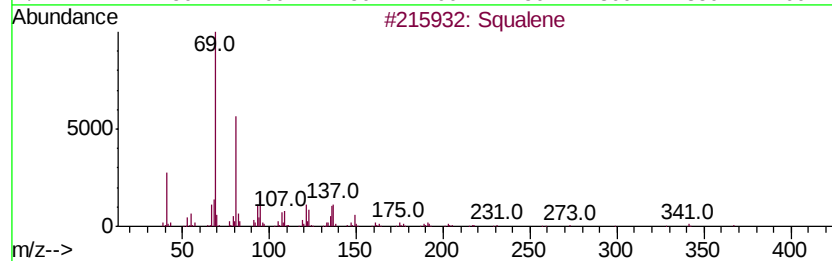
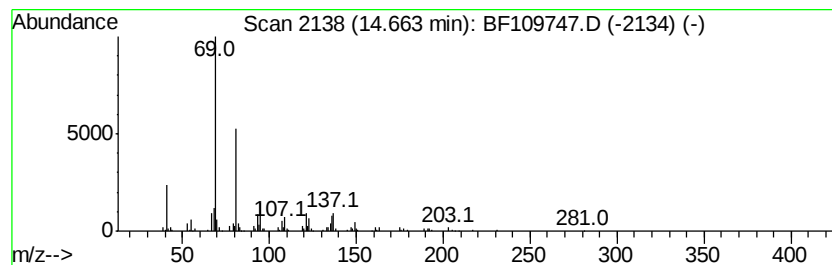
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Squalene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.66	5.06 ng	163157	Perylene-d12	15.23

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Squalene	410	C30H50	000111-02-4	91
2			Supraene	410	C30H50	007683-64-9	91
3			1,5,9-Decatriene, 2,3,5,8-tetram...	192	C14H24	230646-72-7	72
4			(.+/-)-Lavandulol, pentafluorop...	300	C13H17F5O2	1000352-63-1	64
5			1,5-Heptadiene, 3,4-dimethyl-	124	C9H16	1000061-77-9	50



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF101018\
Data File : BF109747.D
Acq On : 10 Oct 2018 17:27
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Sample : J5369-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
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ClientSampleId :
CONCRETE-PILE

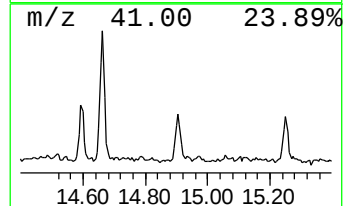
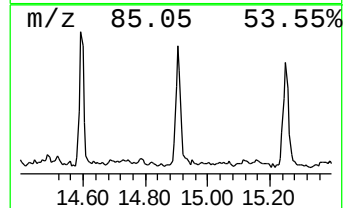
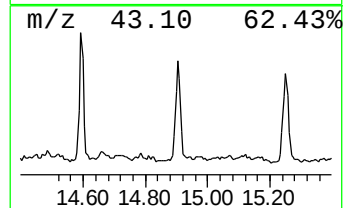
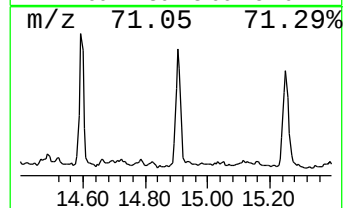
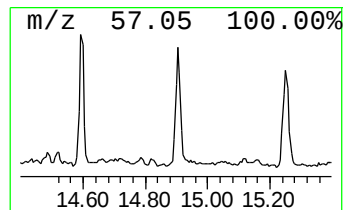
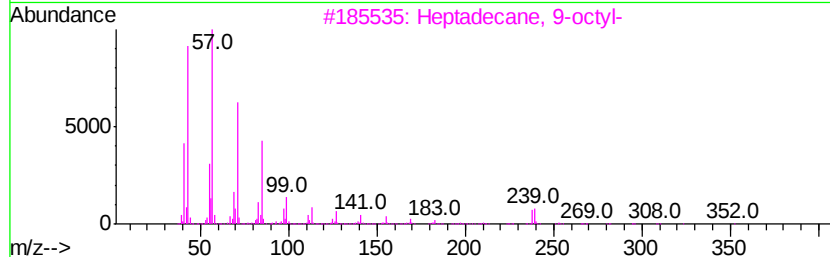
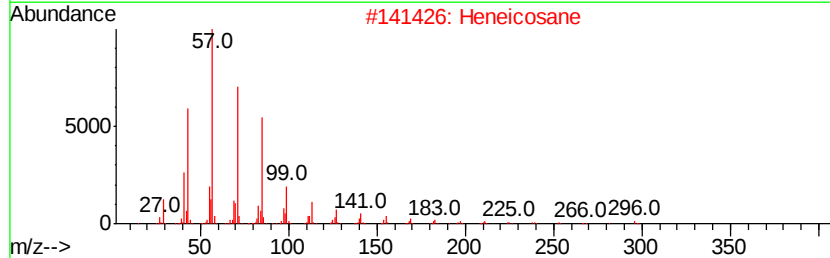
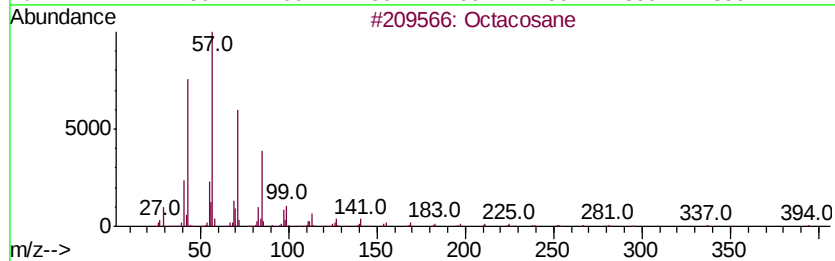
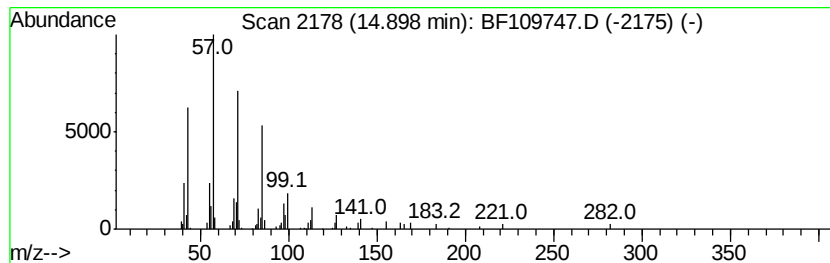
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Octacosane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.90	3.41 ng	109868	Perylene-d12	15.23

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octacosane	394	C28H58	000630-02-4	91
2		Heneicosane	296	C21H44	000629-94-7	91
3		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
4		triacontane	422	C30H62	000638-68-6	91
5		Hexadecane	226	C16H34	000544-76-3	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF101018\
Data File : BF109747.D
Acq On : 10 Oct 2018 17:27
Operator : JU/SJ
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ALS Vial : 11 Sample Multiplier: 1

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ClientSampleId :
CONCRETE-PILE

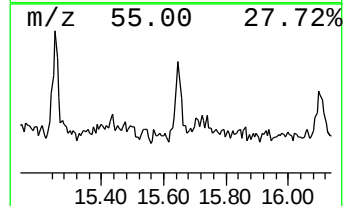
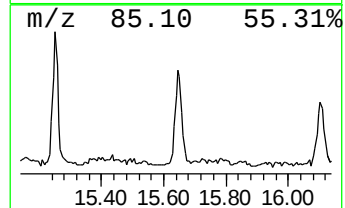
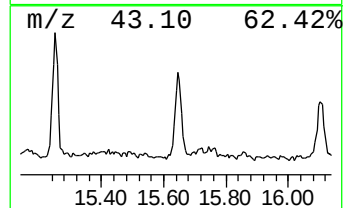
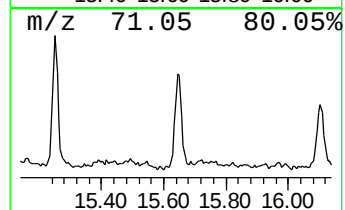
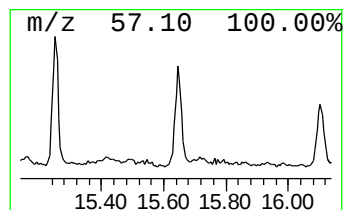
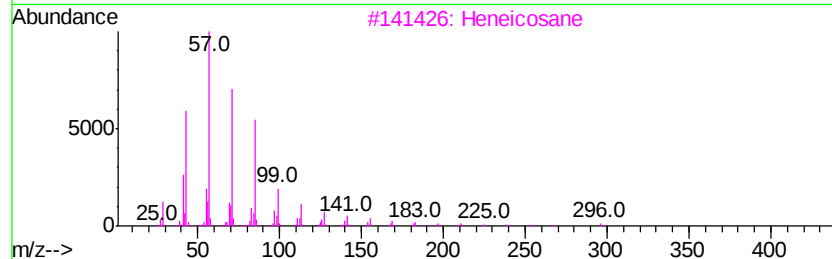
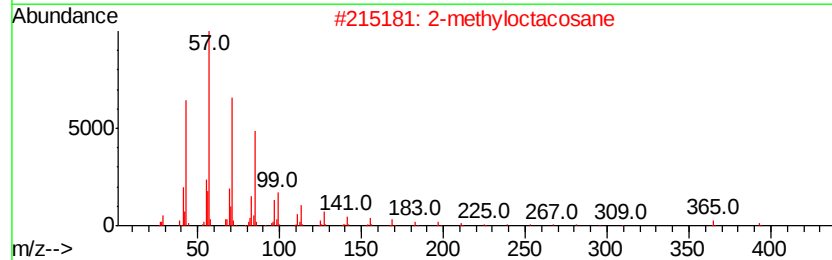
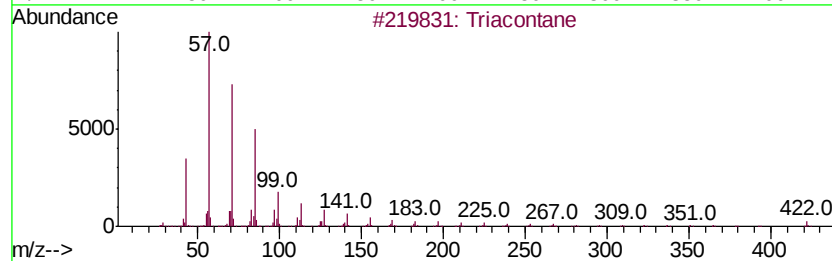
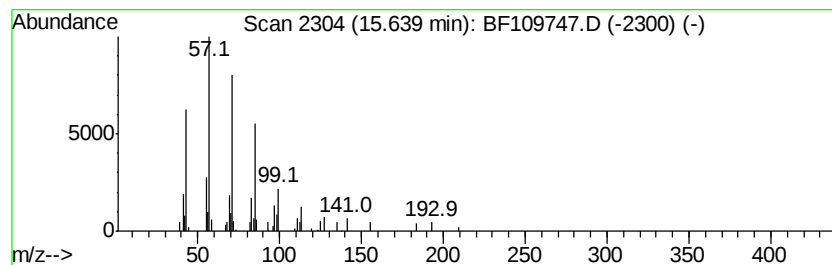
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Docosane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.64	2.88 ng	92805	Perylene-d12	15.23

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Triacontane	422	C30H62	000638-68-6	91
2			2-methyloctacosane	408	C29H60	1000376-72-8	91
3			Heneicosane	296	C21H44	000629-94-7	87
4			Docosane	310	C22H46	000629-97-0	86
5			Hentriacontane	437	C31H64	000630-04-6	81



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF101018\
Data File : BF109747.D
Acq On : 10 Oct 2018 17:27
Operator : JU/SJ
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ALS Vial : 11 Sample Multiplier: 1

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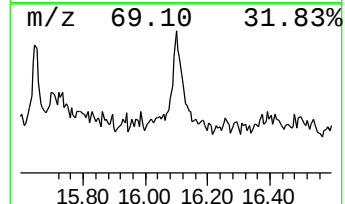
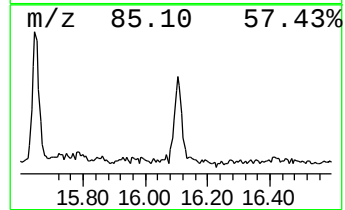
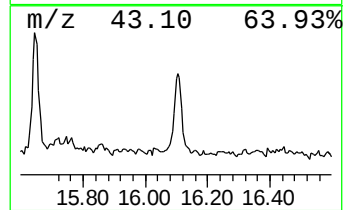
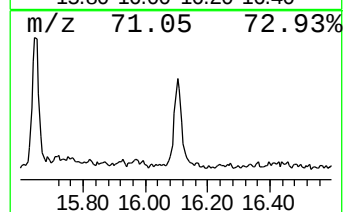
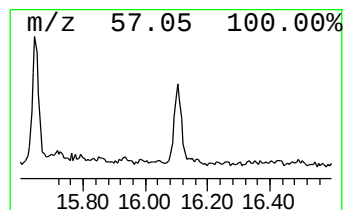
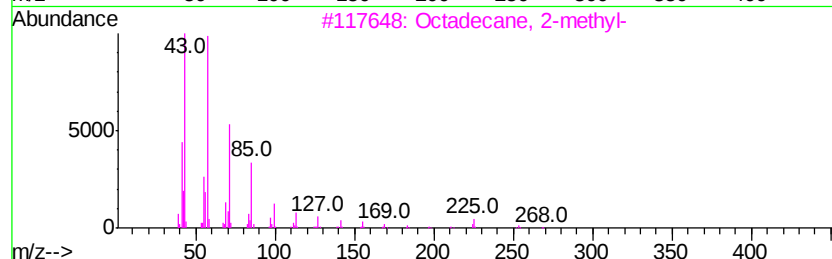
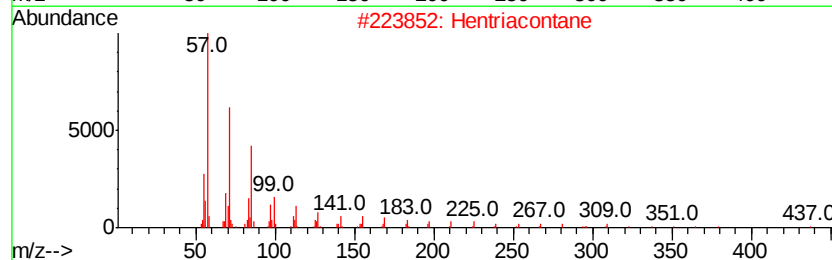
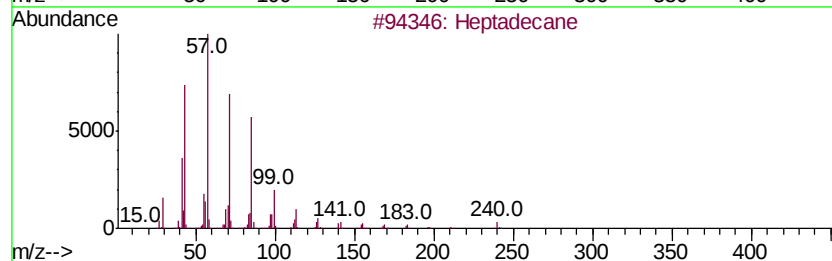
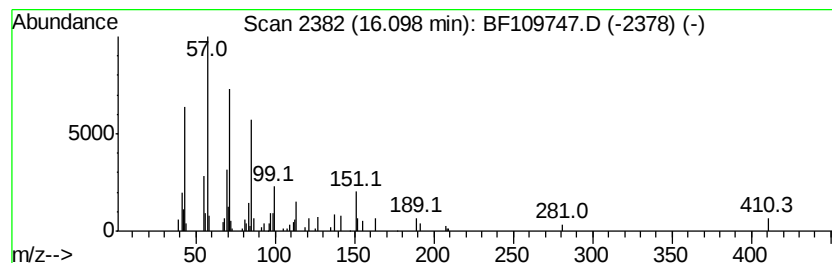
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Hentriacontane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.10	2.41 ng	77729	Perylene-d12	15.23

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heptadecane	240	C17H36	000629-78-7	86
2		Hentriacontane	437	C31H64	000630-04-6	80
3		Octadecane, 2-methyl-	268	C19H40	001560-88-9	78
4		Nonadecane, 2-methyl-	282	C20H42	001560-86-7	78
5		Heneicosane	296	C21H44	000629-94-7	74



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF101018\
 Data File : BF109747.D
 Acq On : 10 Oct 2018 17:27
 Operator : JU/SJ
 Sample : J5369-01
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CONCRETE-PILE

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF100818.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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					#	RT	Resp	Conc
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2-Pentanone, 4-hy...	4.89	64.7	ng	1367750	1	6.70	422924	20.0
unknown6.46	6.46	26.6	ng	561685	1	6.70	422924	20.0
Heptadecane	13.36	2.2	ng	76715	5	13.83	708067	20.0
Octadecane, 1-chl...	13.69	6.3	ng	222474	5	13.83	708067	20.0
Heneicosane	14.00	2.9	ng	101867	5	13.83	708067	20.0
2-methyloctacosane	14.30	3.0	ng	105218	5	13.83	708067	20.0
Triacontane	14.59	3.6	ng	117109	6	15.23	645246	20.0
Squalene	14.66	5.1	ng	163157	6	15.23	645246	20.0
Octacosane	14.90	3.4	ng	109868	6	15.23	645246	20.0
Docosane	15.64	2.9	ng	92805	6	15.23	645246	20.0
Hentriacontane	16.10	2.4	ng	77729	6	15.23	645246	20.0