

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF062918\
 Data File : BF106966.D
 Acq On : 29 Jun 2018 12:13
 Operator : JU/SJ
 Sample : J3709-01
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 EME-FW-TS004-01

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-BF061918.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	496	rBV	121571	142896	9.57%	0.969%
2	5.595	550	554	565	rBV	1266504	1245121	83.37%	8.441%
3	6.595	719	724	737	rVB	1298525	1296770	86.83%	8.792%
4	6.725	742	746	750	rBV	1367341	1316282	88.14%	8.924%
5	6.948	781	784	788	rBV	494140	410978	27.52%	2.786%
6	7.107	807	811	814	rBV	1207414	1027944	68.83%	6.969%
7	7.513	876	880	890	rBV	826981	839020	56.18%	5.688%
8	8.231	998	1002	1016	rBV	527639	535716	35.87%	3.632%
9	9.307	1181	1185	1188	rBV	1736807	1493417	100.00%	10.125%
10	9.701	1248	1252	1257	rBV	183937	167963	11.25%	1.139%
11	9.901	1281	1286	1288	rBV2	33583	40148	2.69%	0.272%
12	9.919	1288	1289	1298	rVV	28530	35505	2.38%	0.241%
13	9.995	1298	1302	1307	rVB	627459	622453	41.68%	4.220%
14	10.401	1368	1371	1374	rVB	46736	33207	2.22%	0.225%
15	10.789	1432	1437	1440	rBV	1013702	996228	66.71%	6.754%
16	10.872	1448	1451	1458	rVB3	38565	41385	2.77%	0.281%
17	11.483	1552	1555	1559	rBV	678564	629782	42.17%	4.270%
18	11.672	1585	1587	1595	rBV4	11986	18499	1.24%	0.125%
19	12.542	1731	1735	1738	rVB4	13397	16768	1.12%	0.114%
20	13.066	1820	1824	1828	rBV	1423091	1291288	86.47%	8.754%
21	13.260	1854	1857	1862	rVB6	15077	24269	1.63%	0.165%
22	13.389	1876	1879	1882	rVB2	32146	29691	1.99%	0.201%
23	13.495	1894	1897	1901	rVB2	63847	57472	3.85%	0.390%
24	13.566	1906	1909	1913	rVB	46863	46257	3.10%	0.314%
25	13.948	1971	1974	1979	rBV	138789	140057	9.38%	0.950%
26	14.142	2003	2007	2014	rBV	578718	567831	38.02%	3.850%
27	14.224	2018	2021	2024	rBV4	23193	22129	1.48%	0.150%
28	14.377	2044	2047	2050	rBV3	30455	32103	2.15%	0.218%
29	14.601	2080	2085	2090	rBV2	66291	98468	6.59%	0.668%
30	14.901	2133	2136	2141	rBV2	52324	66447	4.45%	0.450%
31	14.989	2149	2151	2155	rVB2	26841	25853	1.73%	0.175%
32	15.265	2195	2198	2202	rVB2	94311	98287	6.58%	0.666%
33	15.689	2265	2270	2276	rVB	384664	505365	33.84%	3.426%
34	16.124	2341	2344	2348	rVB2	34859	53050	3.55%	0.360%

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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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35	16.189	2352	2355	2360	rVV4	29748	57369	3.84%	0.389%
36	17.565	2583	2589	2599	rVB5	111560	280954	18.81%	1.905%
37	17.930	2647	2651	2659	rVB9	32107	65494	4.39%	0.444%
38	18.206	2694	2698	2704	rVB7	44794	87959	5.89%	0.596%
39	18.459	2735	2741	2757	rVB7	57492	225876	15.12%	1.531%
40	18.948	2820	2824	2828	rVB6	37269	63870	4.28%	0.433%

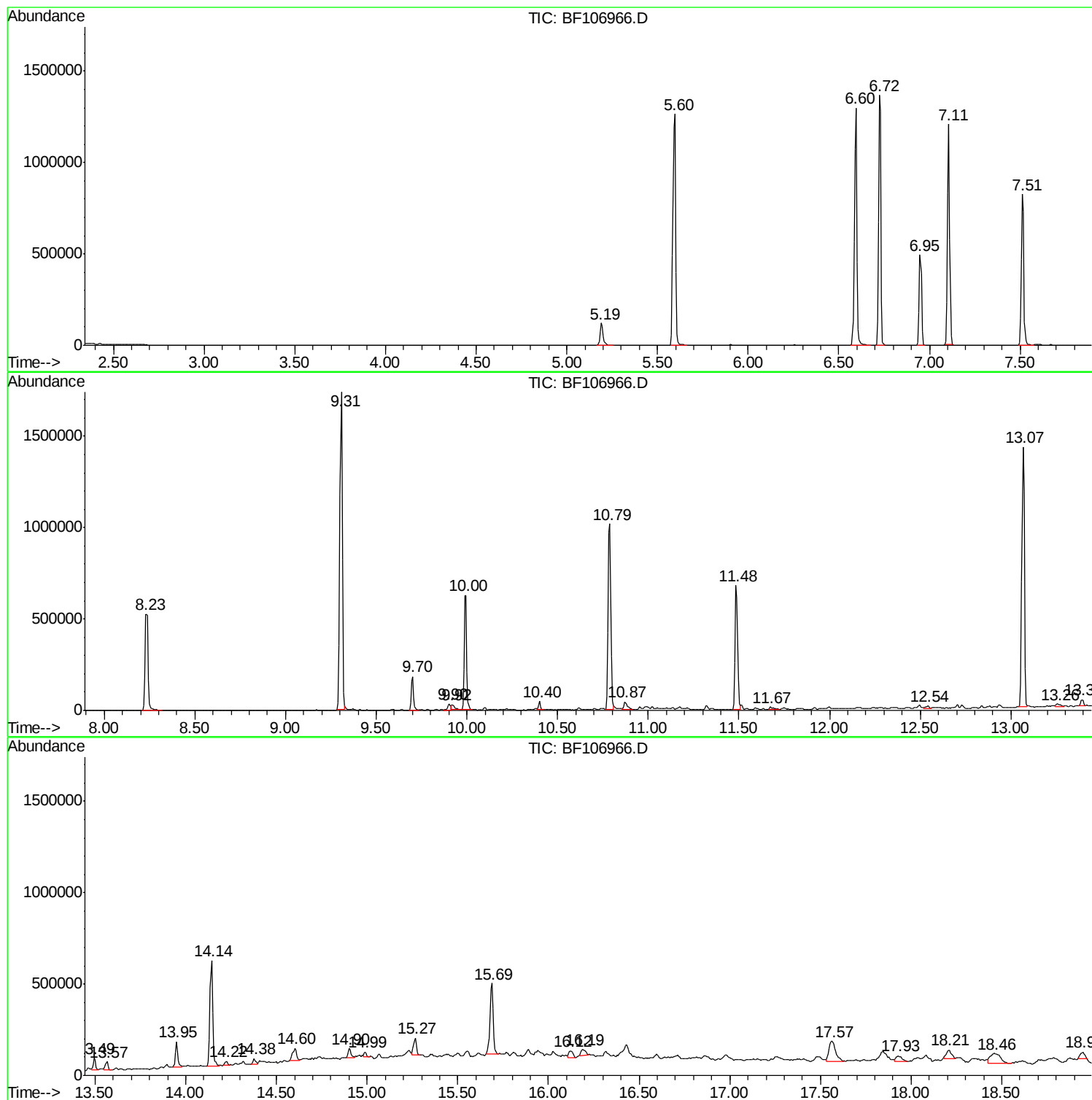
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061918.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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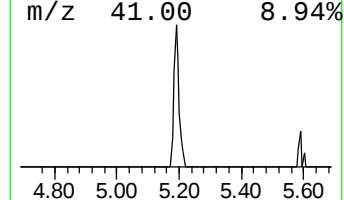
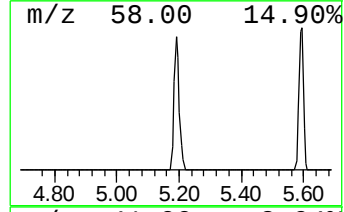
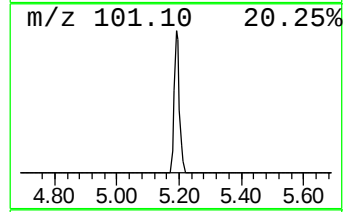
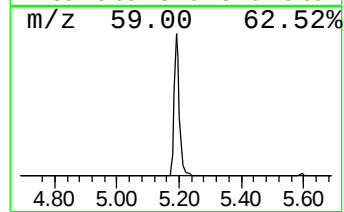
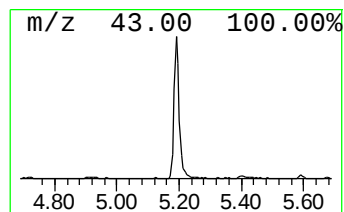
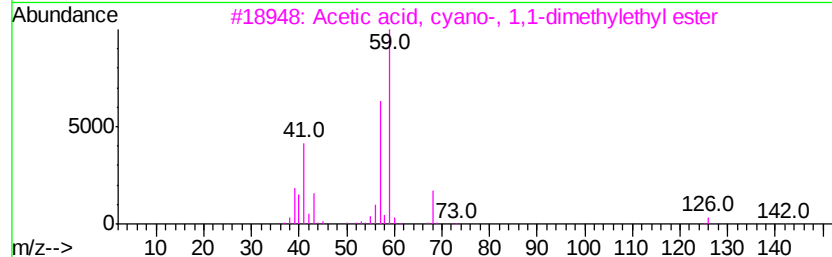
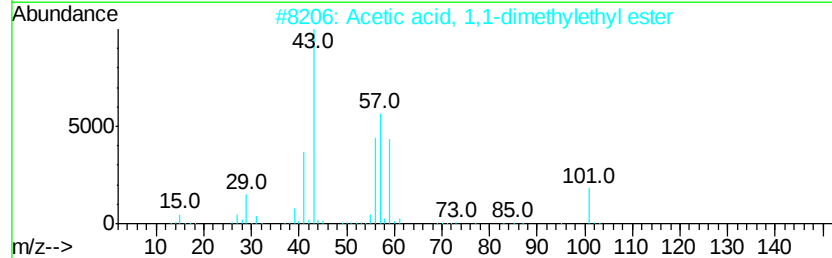
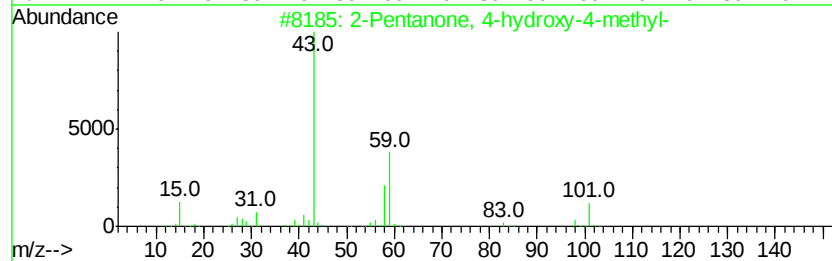
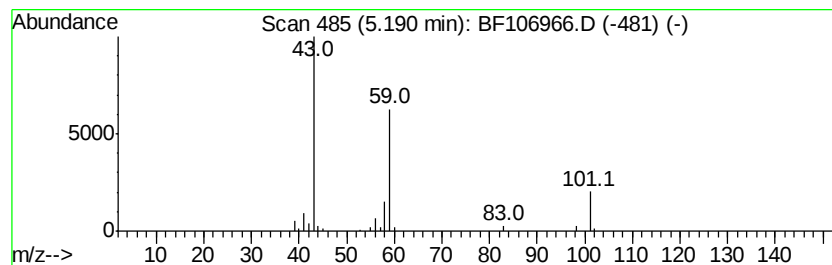
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061918.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.19	6.95 ng	142896	1,4-Dichlorobenzene-d4	6.95

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	40
2			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	39
3			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	23
4			(+)-4-Amino-4,5-dihydro-2(3H)-f...	101	C4H7NO2	016504-58-8	9
5			5-Hexen-2-one	98	C6H10O	000109-49-9	9



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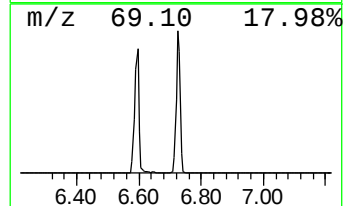
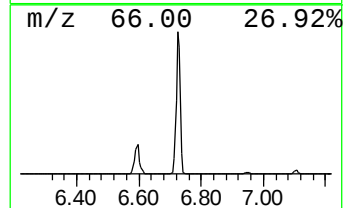
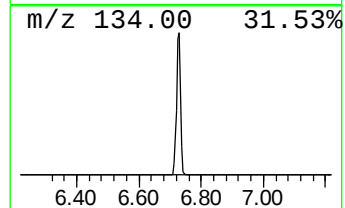
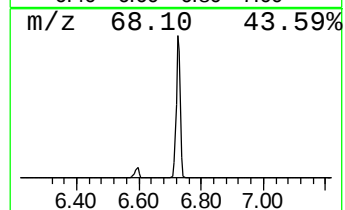
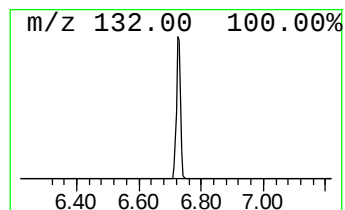
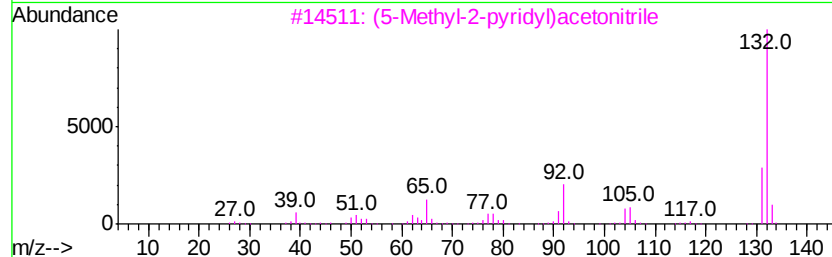
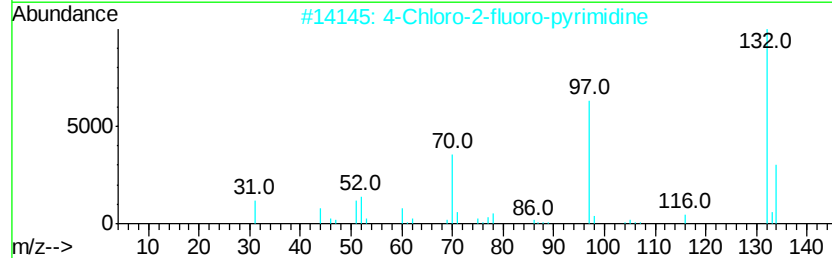
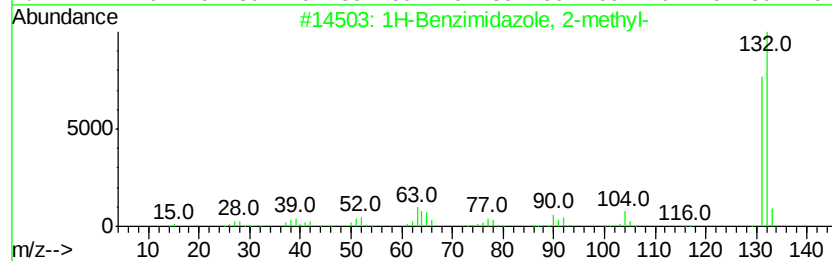
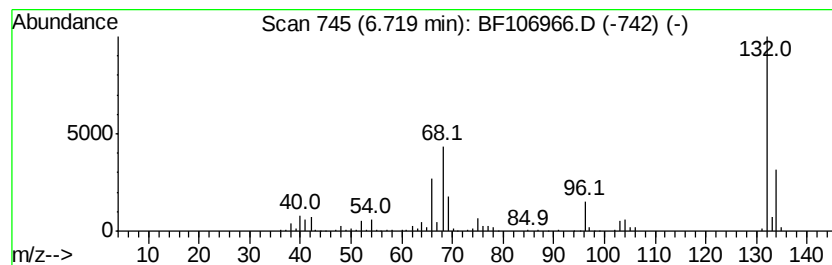
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Peak Number 2 unknown6.72 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.72	64.06 ng	1316280	1,4-Dichlorobenzene-d4	6.95

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	10
2		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9
3		(5-Methyl-2-pyridyl)acetonitrile	132	C8H8N2	1000241-93-9	9
4		2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9
5		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	9



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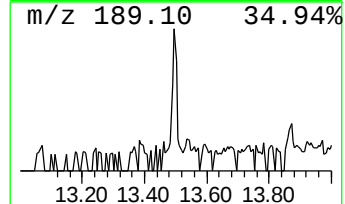
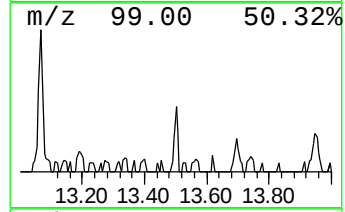
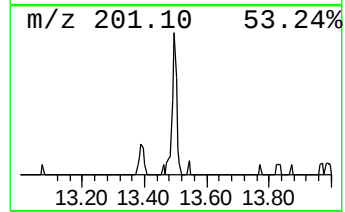
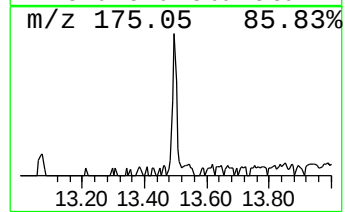
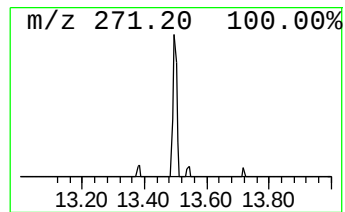
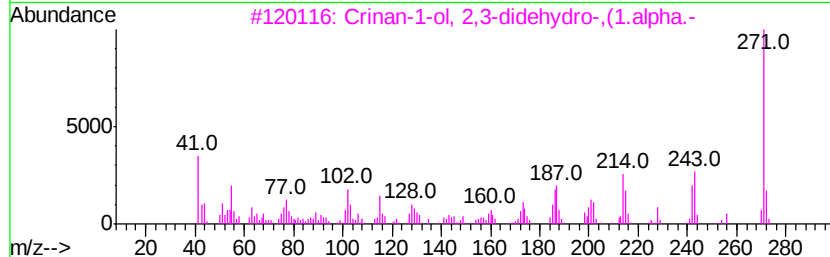
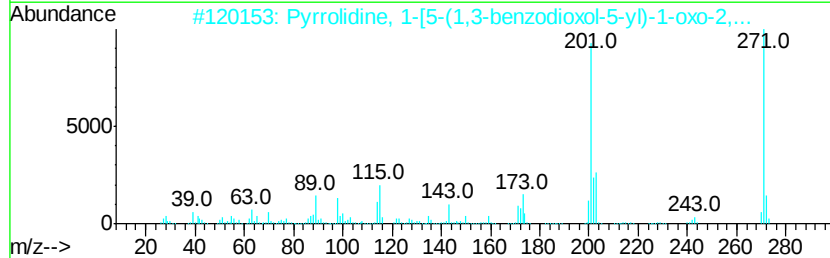
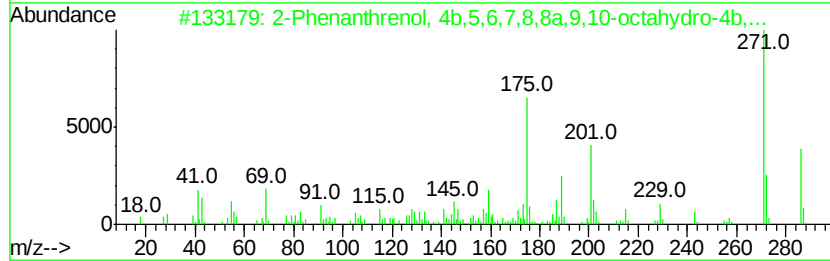
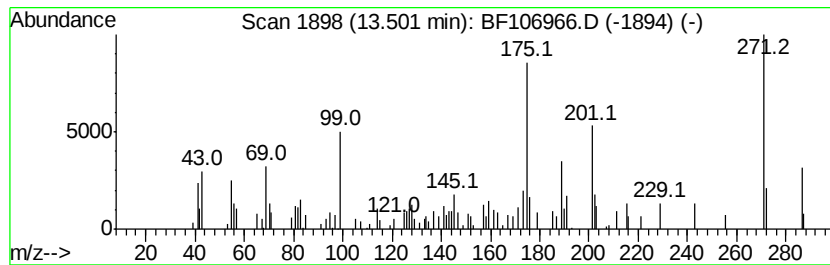
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 4 2-Phenanthrenol, 4b,5,6,7,8... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.		
13.50	2.02 ng	57472	Chrysene-d12	14.14		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Phenanthrenol, 4b,5,6,7,8,8a,9...	286	C20H30O	000511-15-9	70	
2	Pyrrolidine, 1-[5-(1,3-benzodiox...	271	C16H17NO3	025924-78-1	45	
3	Crinan-1-ol, 2,3-didehvdro-, (1.a...	271	C16H17NO3	1000111-31-3	25	
4	Dihydroapohemanthamine	271	C16H17NO3	001153-01-1	25	
5	Dihydronormorphinone	271	C16H17NO3	014696-23-2	18	



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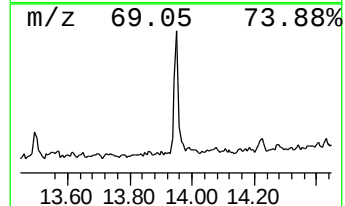
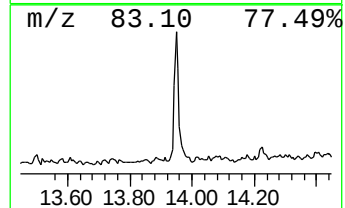
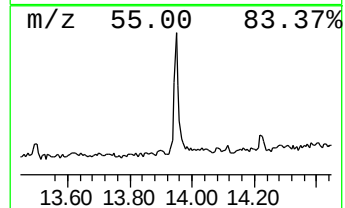
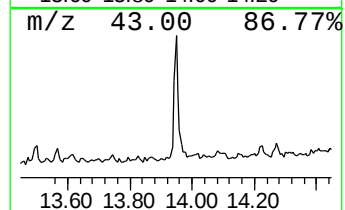
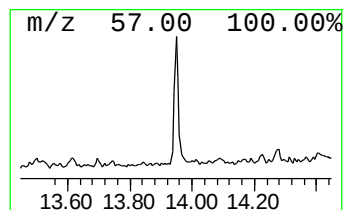
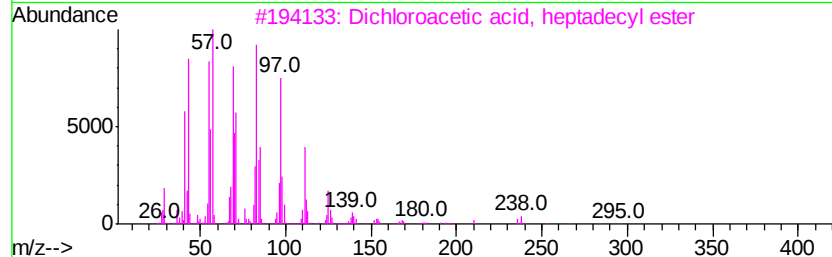
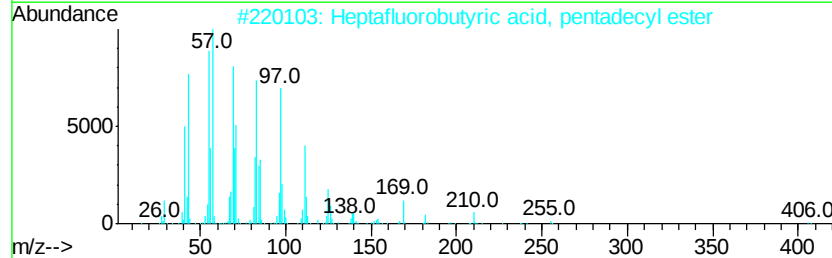
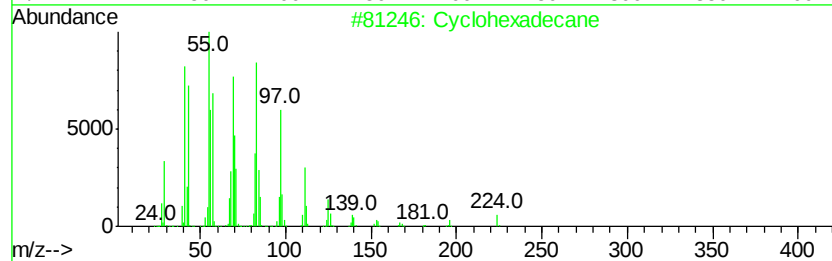
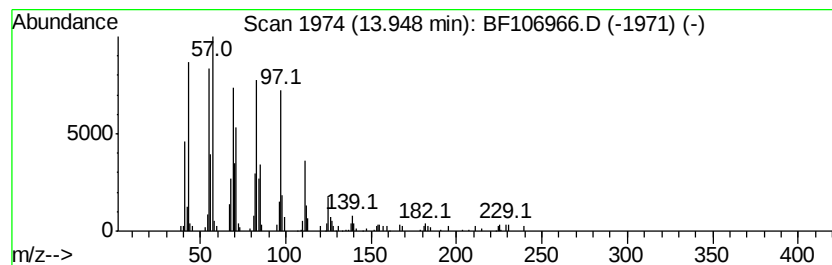
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TIC Integration Parameters: LSCINT.P

Peak Number 5 Cyclohexadecane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.95	4.93 ng	140057	Chrysene-d12	14.14

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cyclohexadecane	224	C16H32	000295-65-8	98
2		Heptafluorobutyric acid, pentade...	424	C19H31F7O2	959261-23-5	95
3		Dichloroacetic acid, heptadecyl ...	366	C19H36Cl2O2	1000282-98-2	95
4		Pentafluoropropionic acid, penta...	374	C18H31F5O2	959092-08-1	95
5		Trifluoroacetic acid, pentadecyl...	324	C17H31F3O2	959010-23-2	95



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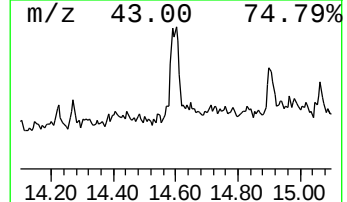
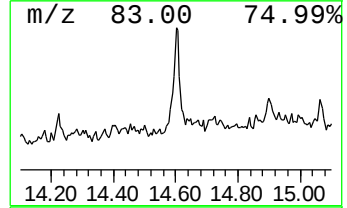
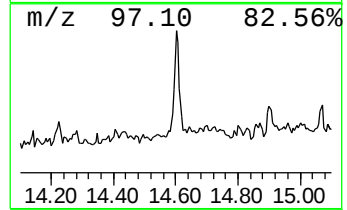
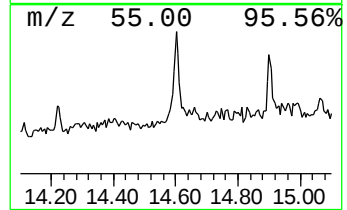
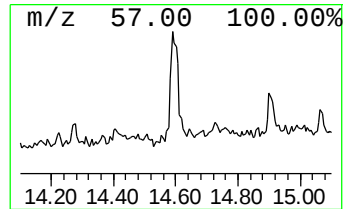
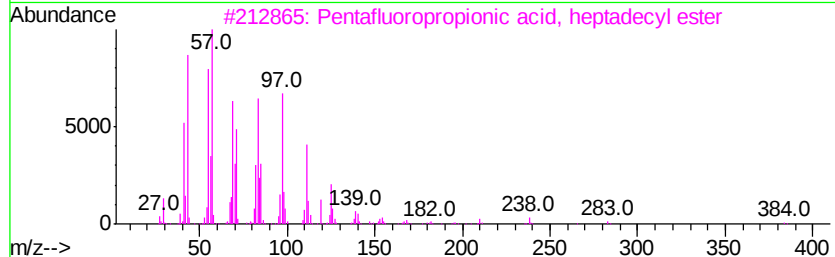
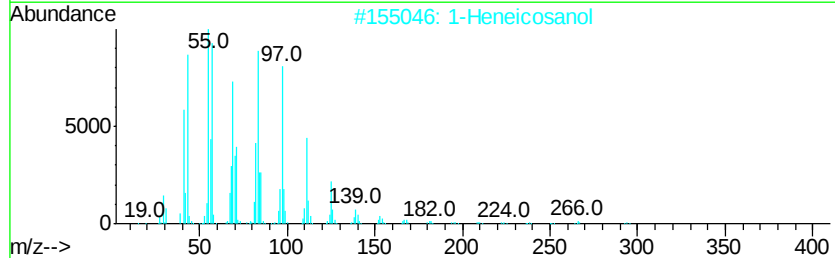
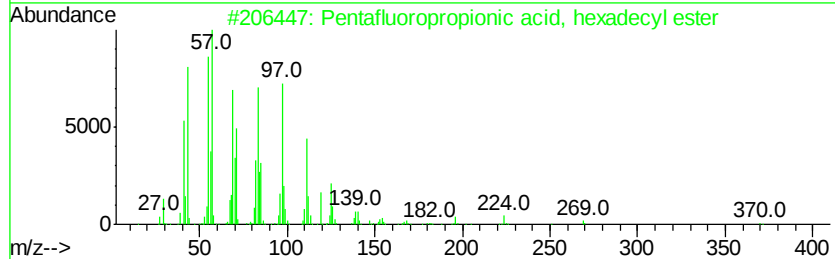
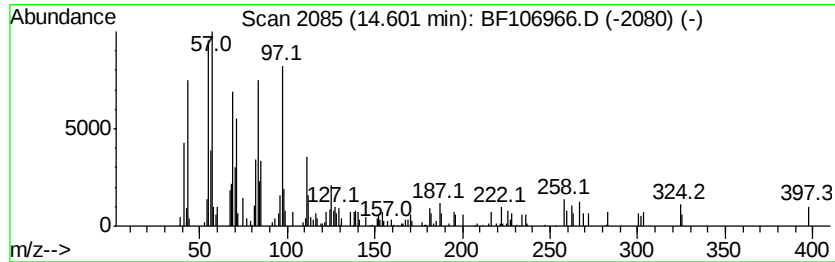
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Pentafluoropropionic acid, ... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.60	3.47 ng	98468	Chrysene-d12	14.14

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentafluoropropionic acid, hexad...	388	C19H33F5O2	006222-07-7	93
2		1-Heneicosanol	312	C21H44O	015594-90-8	93
3		Pentafluoropropionic acid, hepta...	402	C20H35F5O2	959218-78-1	93
4		n-Tetracosanol-1	354	C24H50O	000506-51-4	90
5		Heptadecyl trifluoroacetate	352	C19H35F3O2	1000351-87-0	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062918\
Data File : BF106966.D
Acq On : 29 Jun 2018 12:13
Operator : JU/SJ
Sample : J3709-01
Misc :
ALS Vial : 6 Sample Multiplier: 1

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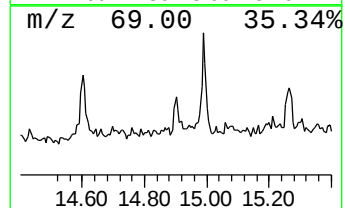
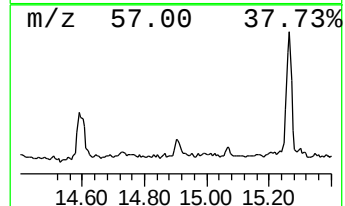
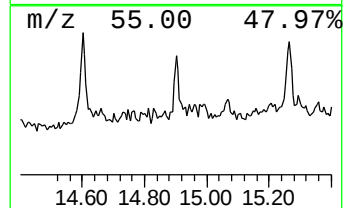
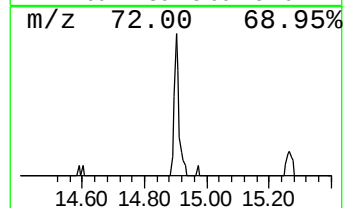
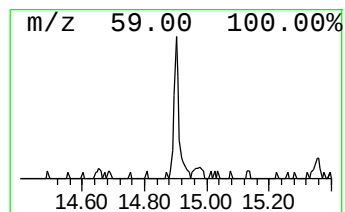
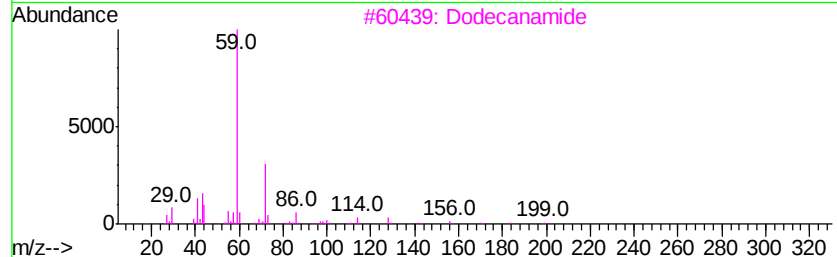
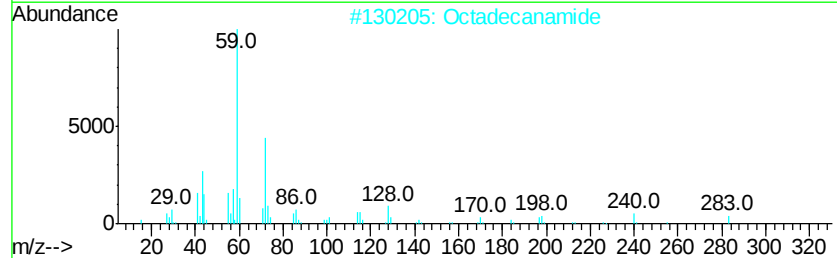
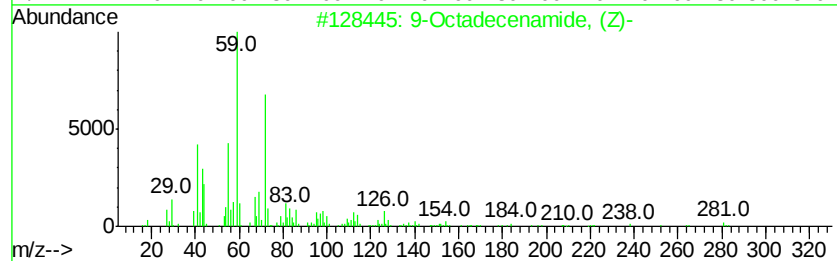
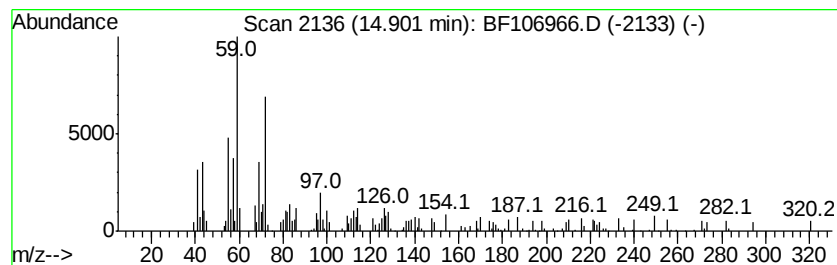
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 9-Octadecenamide, (Z)- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.90	2.34 ng	66447	Chrysene-d12	14.14

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	64
2		Octadecanamide	283	C18H37NO	000124-26-5	59
3		Dodecanamide	199	C12H25NO	001120-16-7	47
4		Diazene, bis[3-(4-methoxycarbonyl...	282	C8H6N6O6	339246-68-3	35
5		Acetic acid, cyanohydroxyimino-,...	128	C4H4N2O3	061295-92-9	35



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062918\
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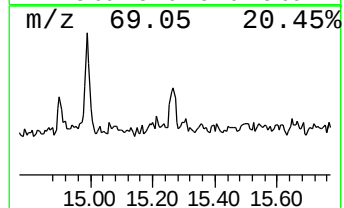
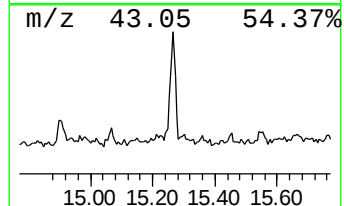
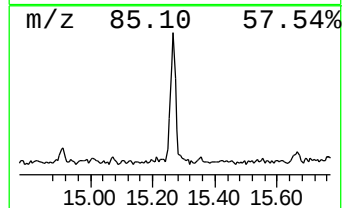
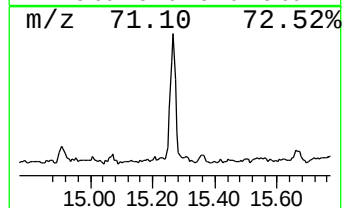
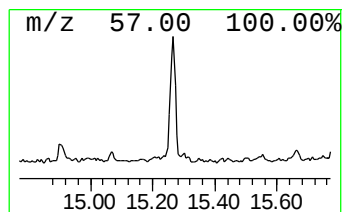
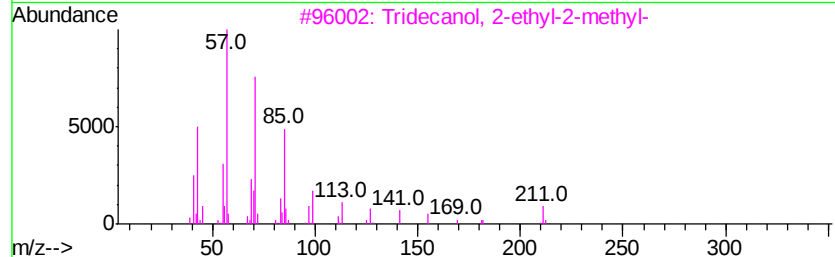
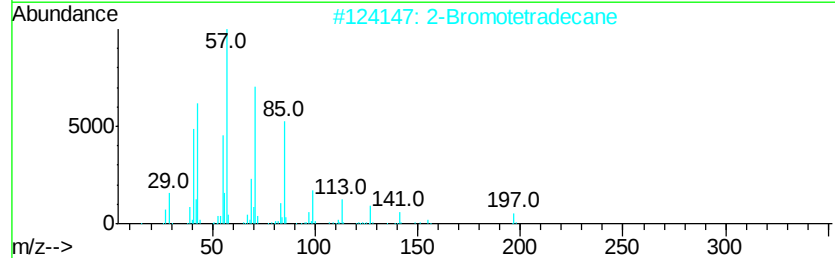
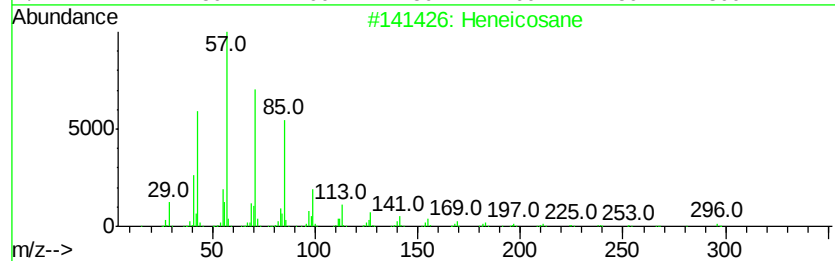
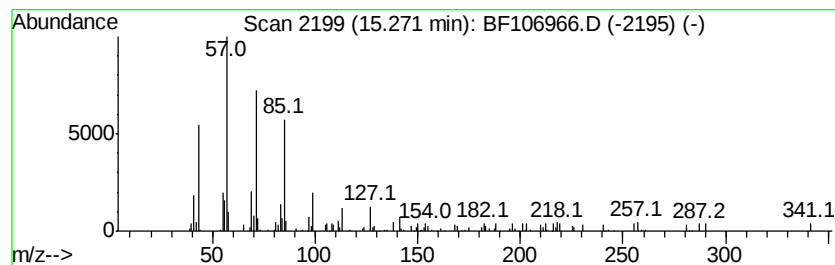
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Heneicosane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.27	3.89 ng	98287	Perylene-d12	15.69

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane		296	C21H44	000629-94-7	91
2	2-Bromotetradecane		276	C14H29Br	074036-95-6	91
3	Tridecanol, 2-ethyl-2-methyl-		242	C16H34O	1000115-66-1	90
4	Octadecane, 1-iodo-		380	C18H37I	000629-93-6	87
5	Octacosane		394	C28H58	000630-02-4	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062918\
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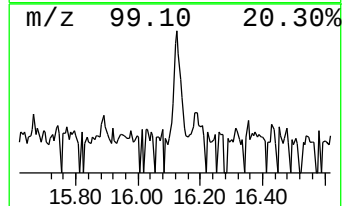
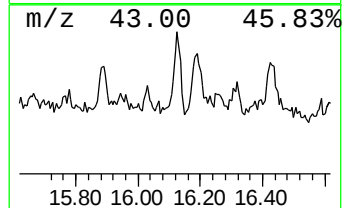
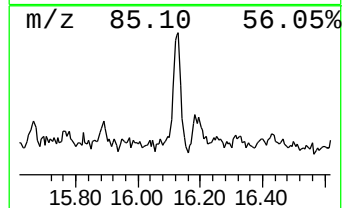
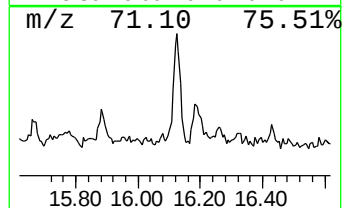
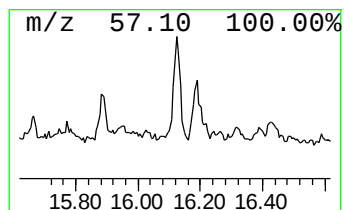
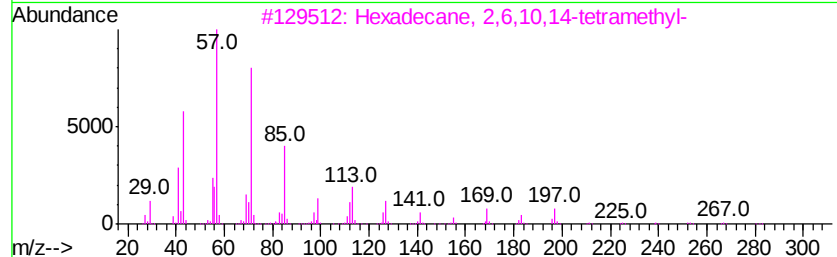
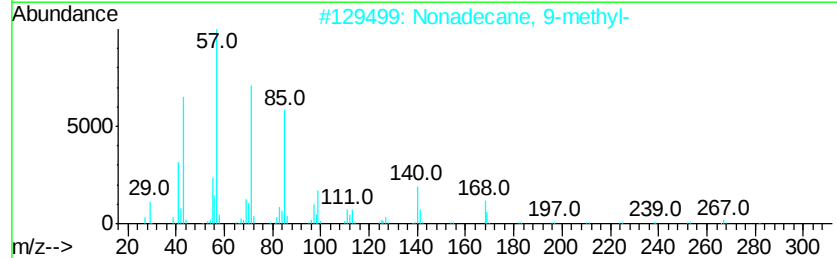
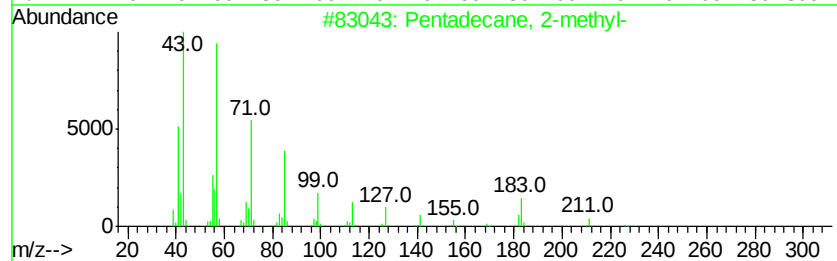
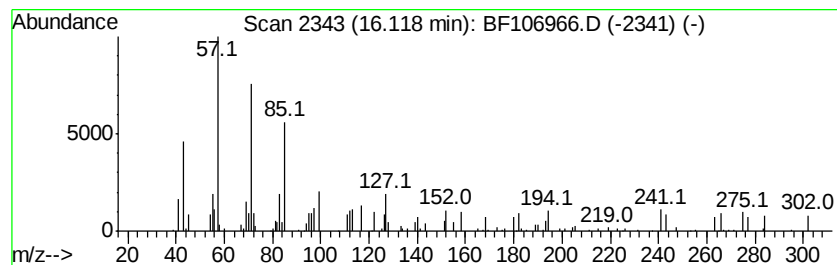
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Pentadecane, 2-methyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.12	2.10 ng	53050	Perylene-d12	15.69

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentadecane, 2-methyl-	226	C16H34	001560-93-6	64
2		Nonadecane, 9-methyl-	282	C20H42	013287-24-6	64
3		Hexadecane, 2,6,10,14-tetramethyl-	282	C20H42	000638-36-8	58
4		Tridecanol, 2-ethyl-2-methyl-	242	C16H34O	1000115-66-1	52
5		Tetracosane	338	C24H50	000646-31-1	52



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062918\
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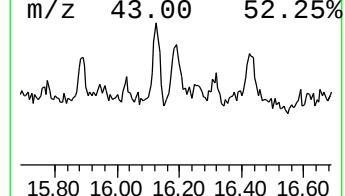
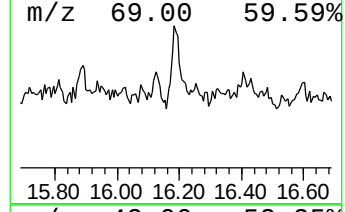
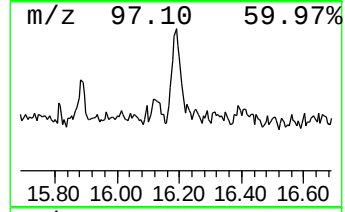
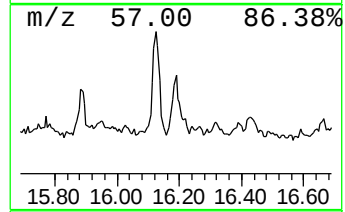
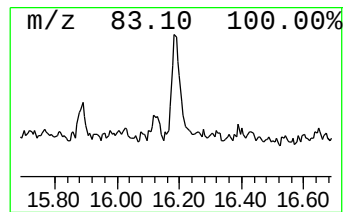
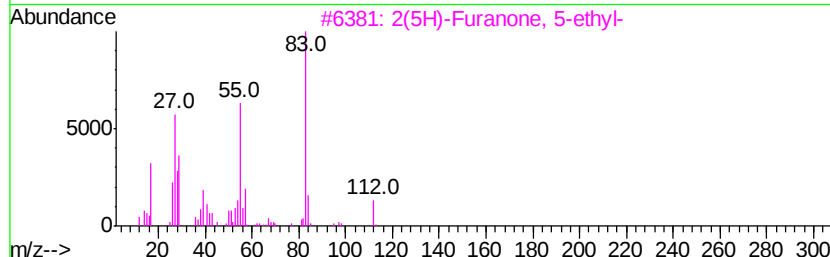
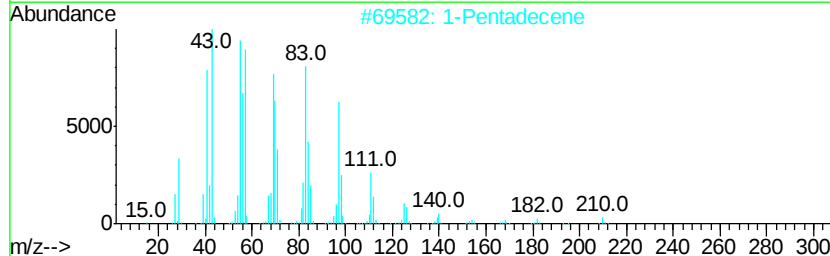
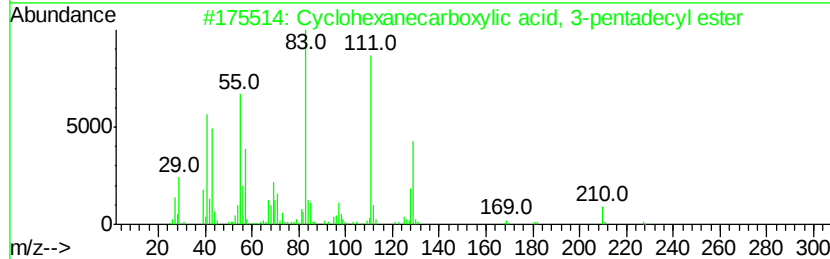
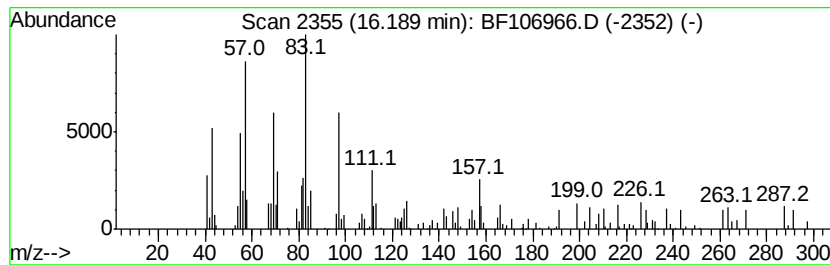
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 unknown16.19 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.19	2.27 ng	57369	Perylene-d12	15.69

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexanecarboxylic acid, 3-pe...	338	C22H42O2	1000280-60-0	38
2			1-Pentadecene	210	C15H30	013360-61-7	38
3			2(5H)-Furanone, 5-ethyl-	112	C6H8O2	002407-43-4	30
4			Cyclohexanol, 1-ethenyl-	126	C8H14O	001940-19-8	30
5			Cyclohexane, iodo-	210	C6H11I	000626-62-0	27



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062918\
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ClientSampleId :
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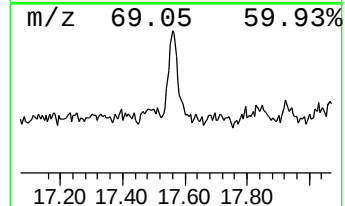
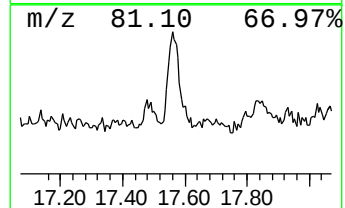
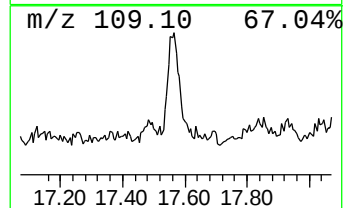
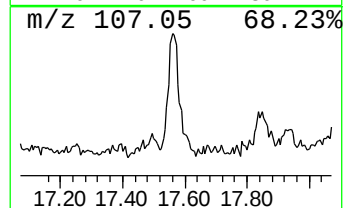
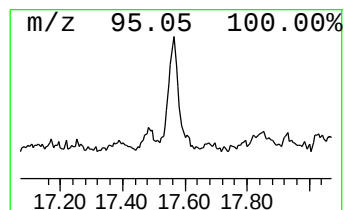
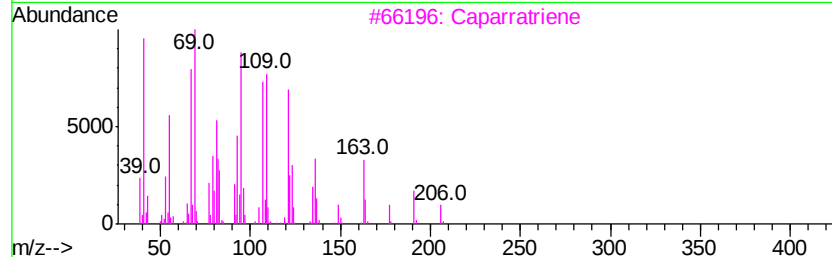
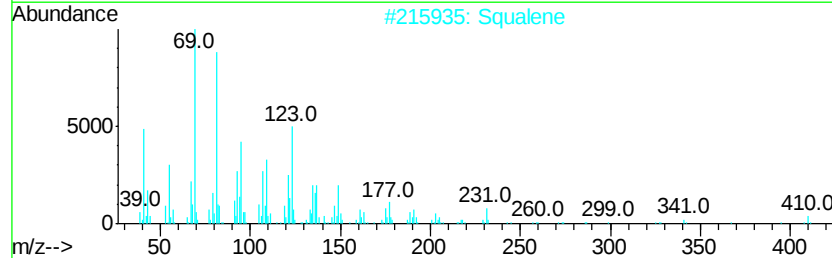
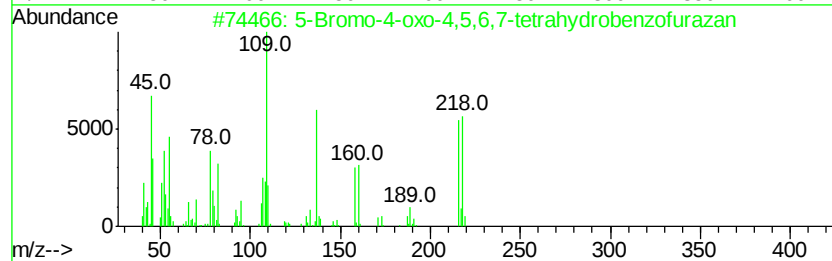
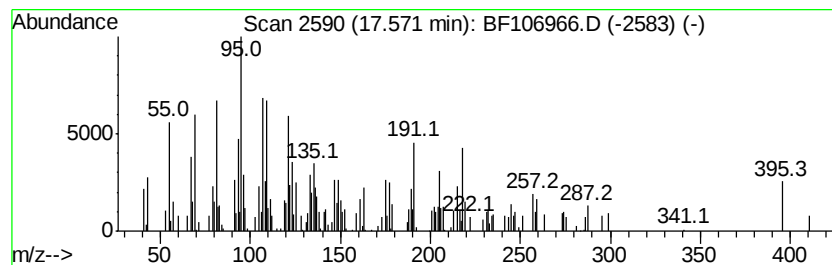
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 5-Bromo-4-oxo-4,5,6,7-tetra... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.57	11.12 ng	280954	Perylene-d12	15.69

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Bromo-4-oxo-4,5,6,7-tetrahydro...	216	C6H5BrN2O2	300574-36-1	92
2		Squalene	410	C30H50	000111-02-4	55
3		Caparratriene	206	C15H26	1000374-08-7	38
4		(2Z,4E)-3,7,11-Trimethyl-2,4,10-...	206	C15H26	172549-29-0	35
5		2,2,6-Trimethyl-1-(2-methyl-cycl...	218	C15H22O	1000188-72-8	35



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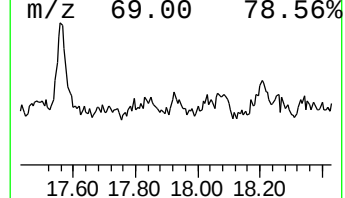
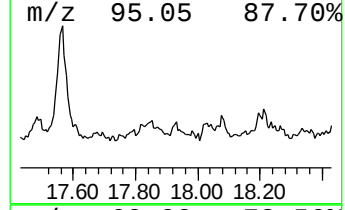
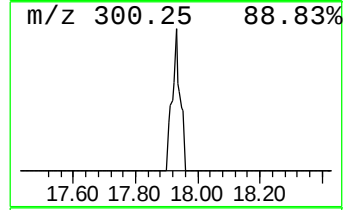
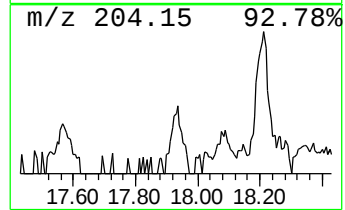
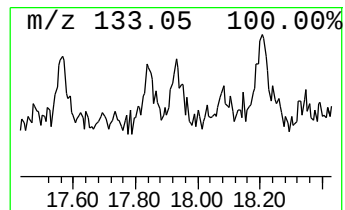
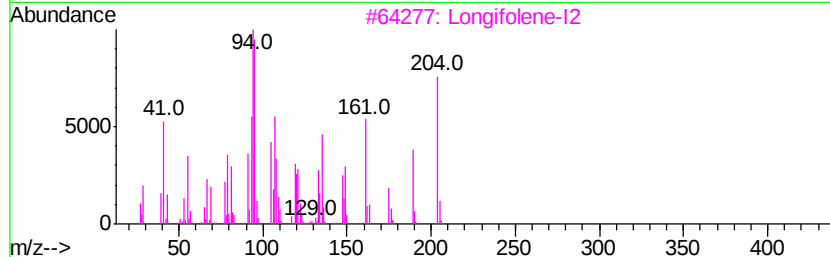
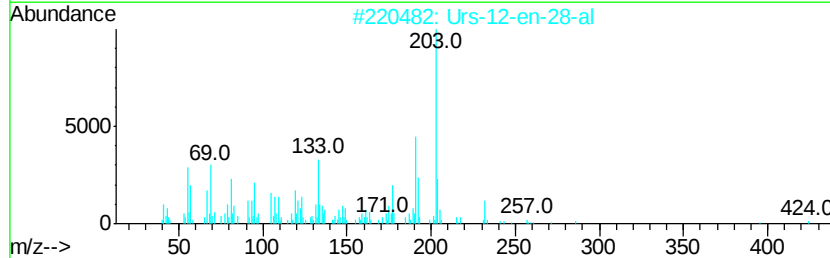
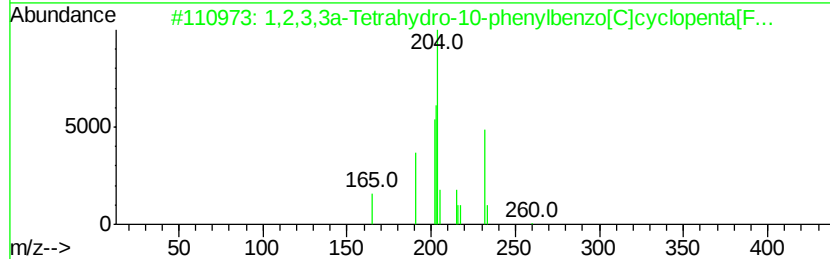
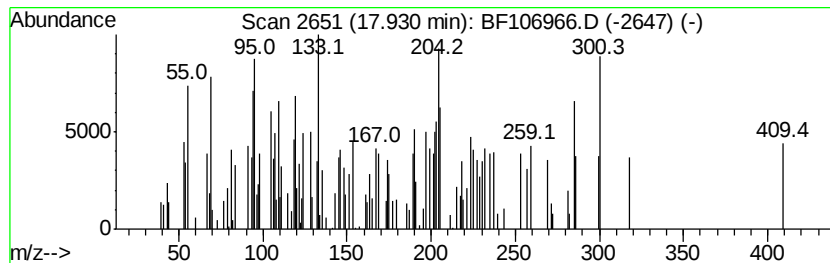
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 12 unknown17.93 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.93	2.59 ng	65494	Perylene-d12	15.69

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,2,3,3a-Tetrahydro-10-phenylben...	260	C18H16N2	028749-06-6	22
2		Urs-12-en-28-al	424	C30H48O	013250-38-9	16
3		Longifolene-I2	204	C15H24	1000162-76-7	10
4		Cyclopropanecarboxamide, 2,2-dib...	337	C11H17Br2NO	1000263-98-9	10
5		4,4'-Bis(tetrahydrothiopyran)	202	C10H18S2	116196-83-9	10



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062918\
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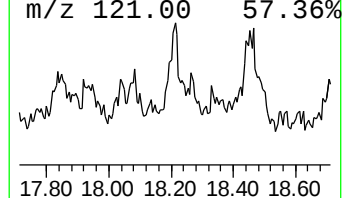
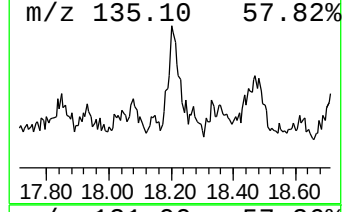
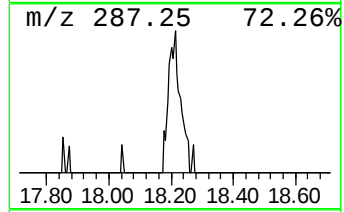
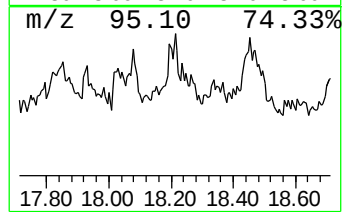
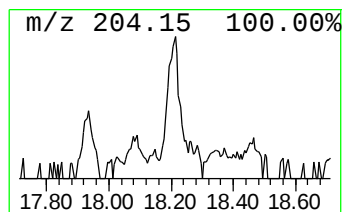
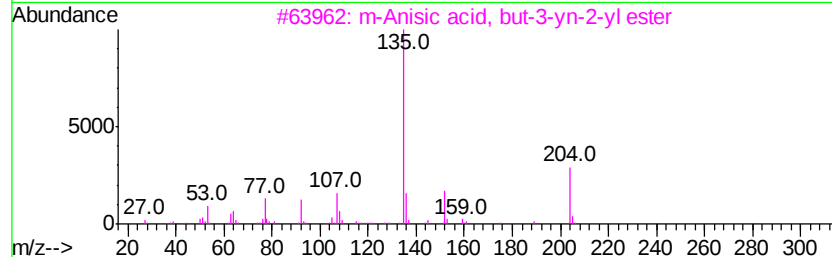
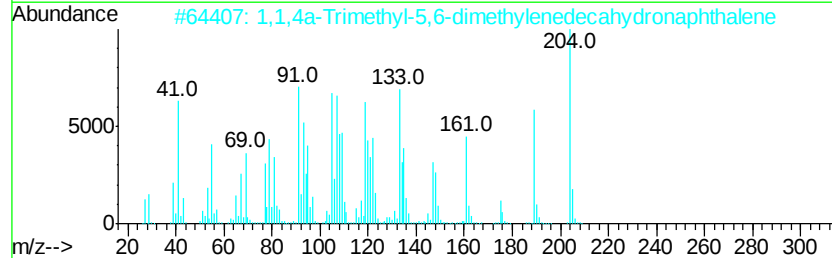
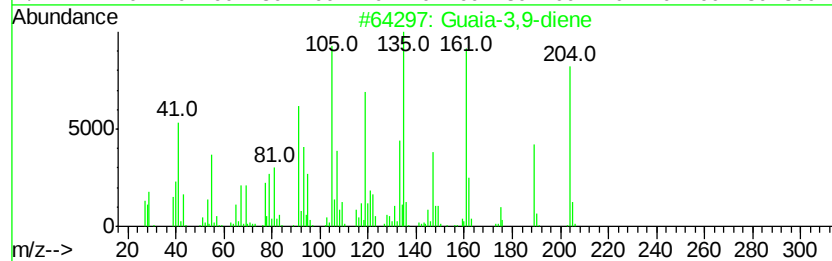
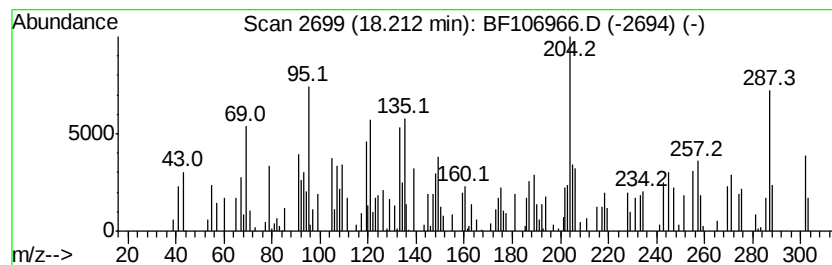
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 unknown18.21 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.21	3.48 ng	87959	Perylene-d12	15.69

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Guaia-3,9-diene	204	C15H24	000489-83-8	46
2		1,1,4a-Trimethyl-5,6-dimethylene...	204	C15H24	1000193-60-8	40
3		m-Anisic acid, but-3-yn-2-yl ester	204	C12H12O3	1000292-59-1	38
4		2H-Cyclopentacyclooctene, 4,5,6,...	204	C15H24	1000221-85-8	38
5		Naphthalene, 1,2,3,5,6,7,8,8a-oc...	204	C15H24	004630-07-3	35



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062918\
Data File : BF106966.D
Acq On : 29 Jun 2018 12:13
Operator : JU/SJ
Sample : J3709-01
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
EME-FW-TS004-01

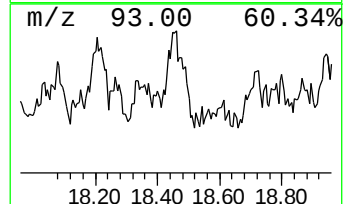
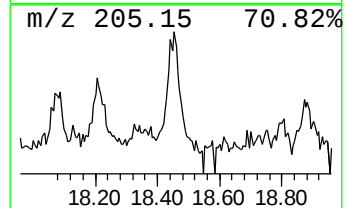
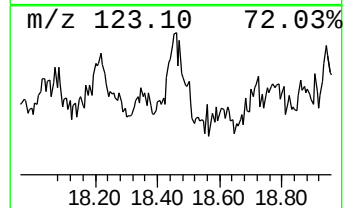
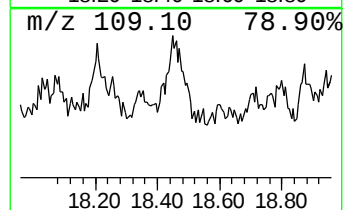
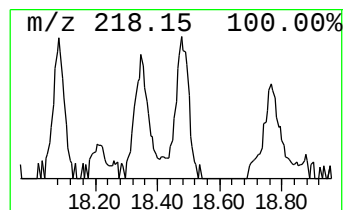
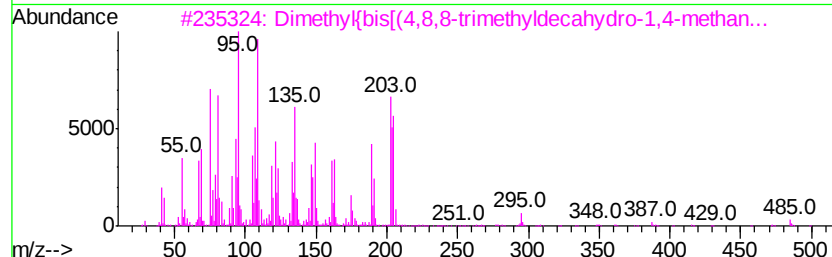
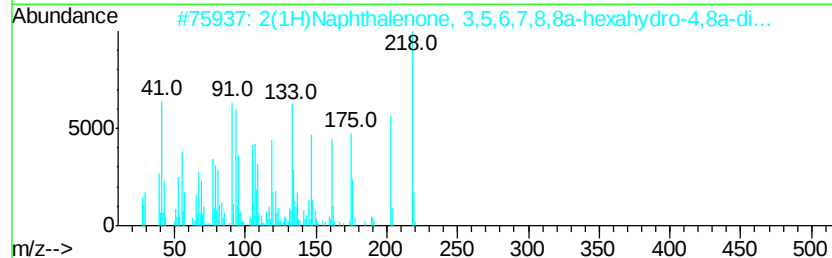
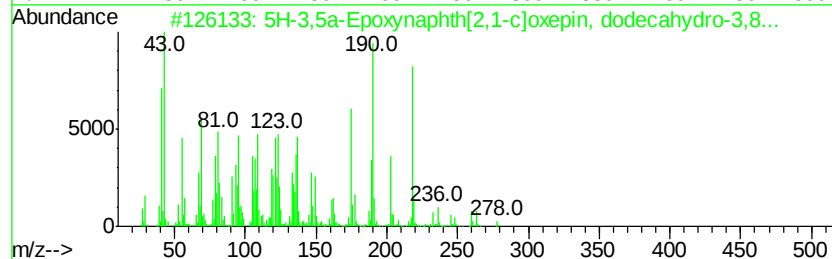
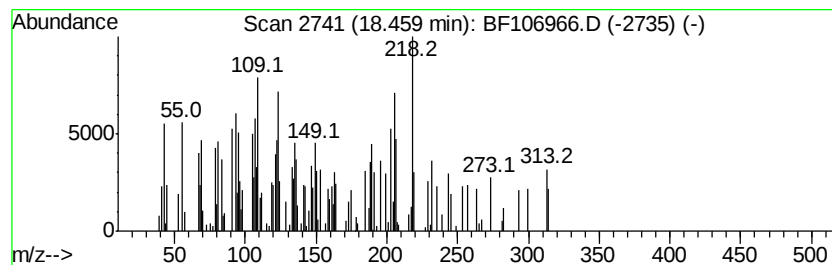
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 unknown18.46 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.46	8.94 ng	225876	Perylene-d12	15.69

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5H-3,5a-Epoxy naphth[2,1-c]oxepin...	278	C18H30O2	001153-34-0	18
2		2(1H)Naphthalenone, 3,5,6,7,8,8a...	218	C15H22O	1000188-66-5	15
3		Dimethylbis[4,8,8-trimethyldec...	500	C32H56O2Si	1000364-69-0	14
4		2,4,4-Trimethyl-3-(3-methylbuta...	206	C14H22O	097452-05-6	14
5		6-Isopropenyl-4,8a-dimethyl-4a,5...	218	C15H22O	086917-79-5	11



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062918\
Data File : BF106966.D
Acq On : 29 Jun 2018 12:13
Operator : JU/SJ
Sample : J3709-01
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
EME-FW-TS004-01

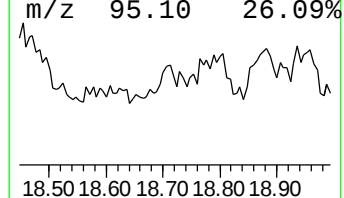
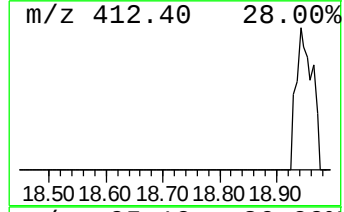
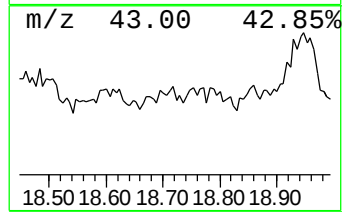
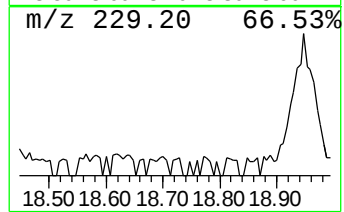
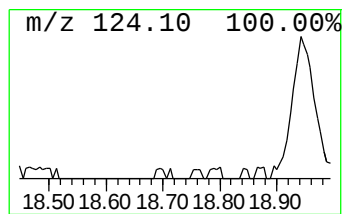
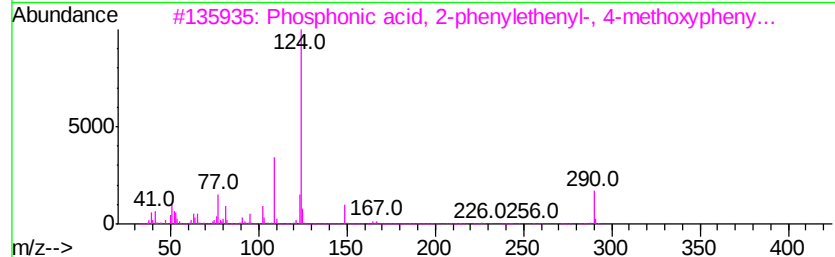
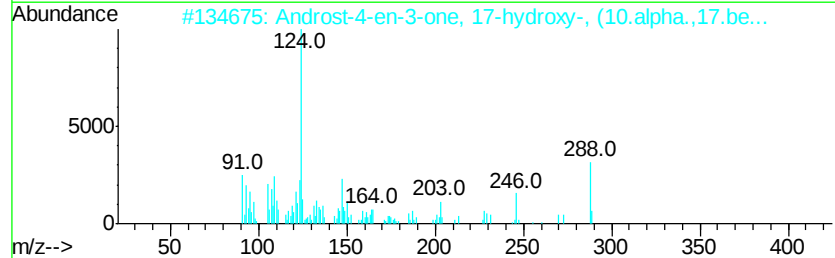
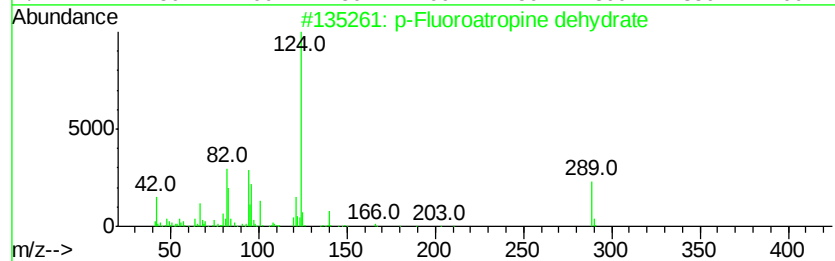
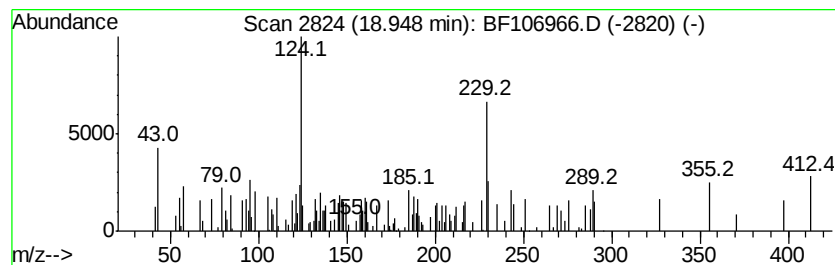
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 unknown18.95 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.95	2.53 ng	63870	Perylene-d12	15.69

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	p-Fluoroatropine dehydrate	289	C17H20FN02	1000212-74-8	27
2		Androst-4-en-3-one, 17-hydroxy-,...	288	C19H28O2	000604-39-7	25
3		Phosphonic acid, 2-phenylethenvl...	290	C15H15O4P	330855-58-8	22
4		Testosterone	288	C19H28O2	000058-22-0	22
5		Pyrrolizin-1-one, 7(S)-[(2-ethyl...	257	C12H19NOS2	1000125-98-4	18



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF062918\
 Data File : BF106966.D
 Acq On : 29 Jun 2018 12:13
 Operator : JU/SJ
 Sample : J3709-01
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 EME-FW-TS004-01

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF061918.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
2-Pentanone, 4-hy...	5.19	7.0	ng	142896	1	6.95	410978	20.0
unknown6.72	6.72	64.1	ng	1316280	1	6.95	410978	20.0
2-Phenanthrenol, ...	13.50	2.0	ng	57472	5	14.14	567831	20.0
Cyclohexadecane	13.95	4.9	ng	140057	5	14.14	567831	20.0
Pentafluoropropio...	14.60	3.5	ng	98468	5	14.14	567831	20.0
9-Octadecenamide, ...	14.90	2.3	ng	66447	5	14.14	567831	20.0
Heneicosane	15.27	3.9	ng	98287	6	15.69	505365	20.0
Pentadecane, 2-me...	16.12	2.1	ng	53050	6	15.69	505365	20.0
unknown16.19	16.19	2.3	ng	57369	6	15.69	505365	20.0
5-Bromo-4-oxo-4,5...	17.57	11.1	ng	280954	6	15.69	505365	20.0
unknown17.93	17.93	2.6	ng	65494	6	15.69	505365	20.0
unknown18.21	18.21	3.5	ng	87959	6	15.69	505365	20.0
unknown18.46	18.46	8.9	ng	225876	6	15.69	505365	20.0
unknown18.95	18.95	2.5	ng	63870	6	15.69	505365	20.0