

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF110118\
 Data File : BF110370.D
 Acq On : 1 Nov 2018 15:55
 Operator : JU/SJ
 Sample : J5741-55
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-14

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-BF102318.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	2.316	31	39	46	rVB	436580	608053	17.28%	2.189%
2	3.599	255	257	267	rBV	56979	90916	2.58%	0.327%
3	5.210	528	531	543	rVB	96792	137370	3.90%	0.495%
4	5.587	590	595	598	rBV	2680329	2779726	79.01%	10.007%
5	6.587	758	765	769	rBV	2610438	3049660	86.68%	10.979%
6	6.728	784	789	792	rBV	2961462	2997568	85.20%	10.791%
7	6.951	824	827	831	rBV	595111	564116	16.03%	2.031%
8	7.110	850	854	858	rBV	2455546	2297447	65.30%	8.271%
9	7.522	918	924	927	rBV	1774402	1888199	53.67%	6.798%
10	8.240	1041	1046	1055	rBV	733272	736265	20.93%	2.651%
11	9.316	1222	1229	1232	rBV	3386746	3518285	100.00%	12.666%
12	9.704	1290	1295	1305	rVB	407650	396476	11.27%	1.427%
13	9.998	1340	1345	1353	rBV	931767	845802	24.04%	3.045%
14	10.792	1475	1480	1483	rBV	2118758	2032599	57.77%	7.317%
15	11.486	1594	1598	1606	rBV	857550	794328	22.58%	2.860%
16	11.992	1681	1684	1689	rBV	52884	52663	1.50%	0.190%
17	13.074	1863	1868	1871	rBV	2707971	2547883	72.42%	9.173%
18	13.945	2012	2016	2030	rBV	167575	220774	6.28%	0.795%
19	14.133	2044	2048	2053	rBV	537824	489053	13.90%	1.761%
20	14.263	2066	2070	2075	rBV	72166	69420	1.97%	0.250%
21	14.569	2118	2122	2130	rBV2	109967	150012	4.26%	0.540%
22	14.886	2171	2176	2180	rBV	116664	121343	3.45%	0.437%
23	15.239	2231	2236	2239	rBV	146025	158696	4.51%	0.571%
24	15.657	2299	2307	2316	rVB2	393966	655935	18.64%	2.361%
25	16.092	2373	2381	2386	rBV	116639	182594	5.19%	0.657%
26	16.163	2388	2393	2399	rVB6	19363	43158	1.23%	0.155%
27	16.627	2466	2472	2478	rBV	56273	99507	2.83%	0.358%
28	17.245	2573	2577	2586	rVB2	27238	56410	1.60%	0.203%
29	17.357	2590	2596	2605	rBV5	68033	141733	4.03%	0.510%
30	17.804	2666	2672	2681	rBV10	18618	51366	1.46%	0.185%

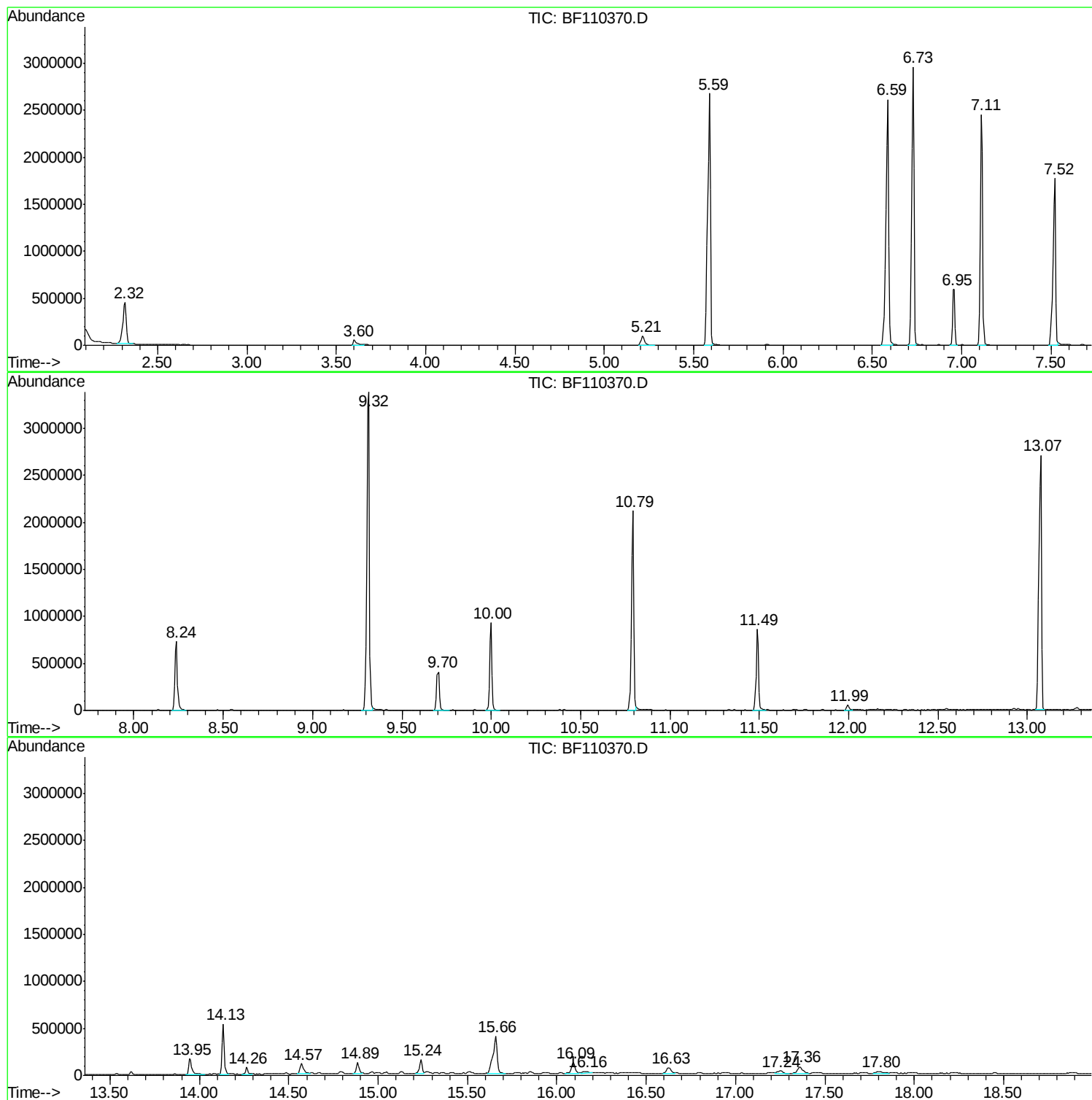
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ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF102318.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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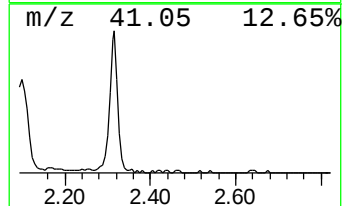
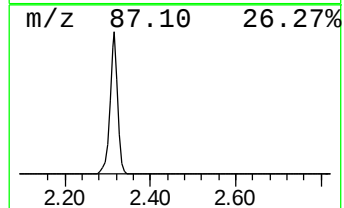
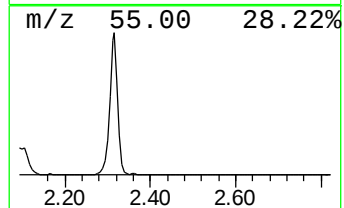
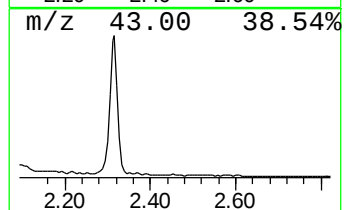
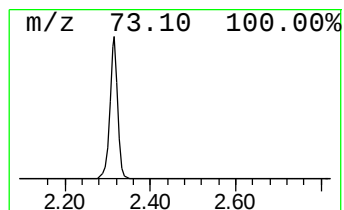
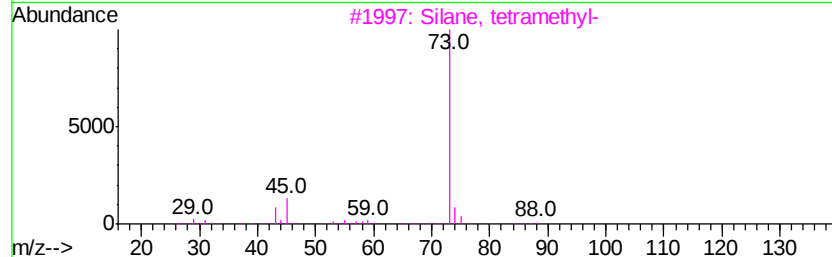
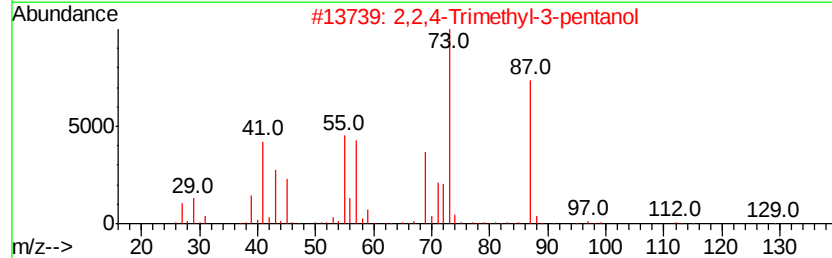
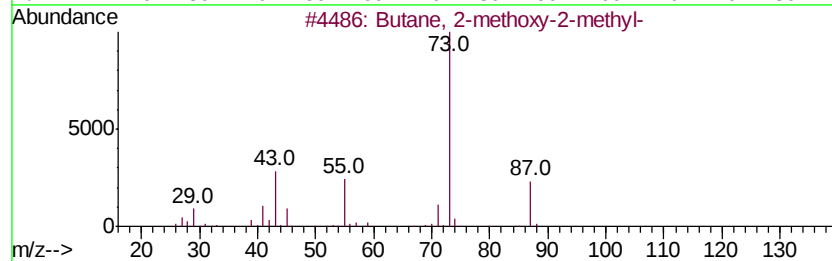
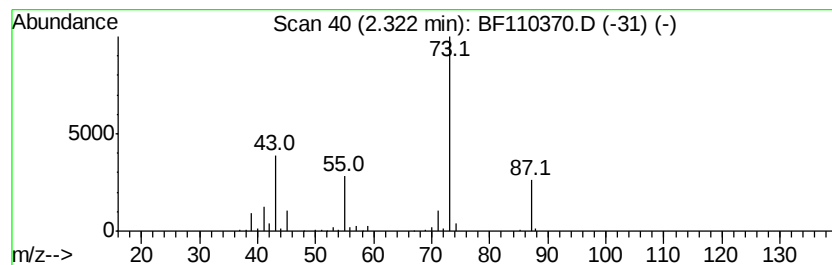
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.32	21.56 ng	608053	1,4-Dichlorobenzene-d4	6.96

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2		2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39
3		Silane, tetramethyl-	88	C4H12Si	000075-76-3	38
4		1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	37
5		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	12



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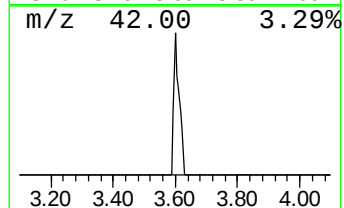
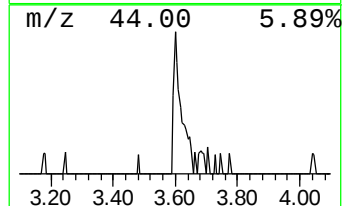
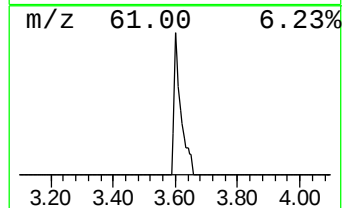
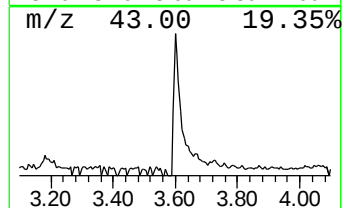
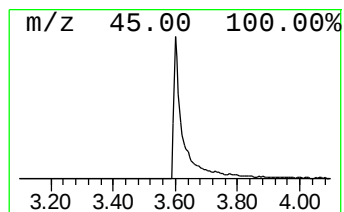
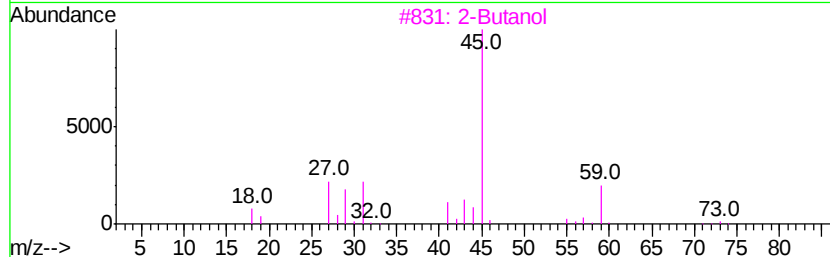
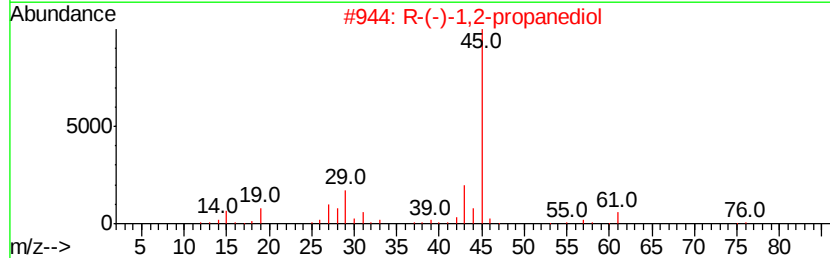
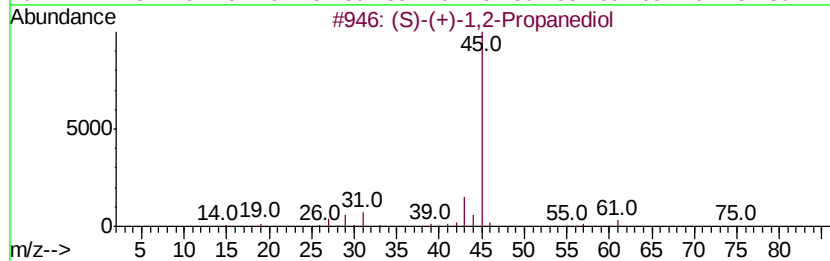
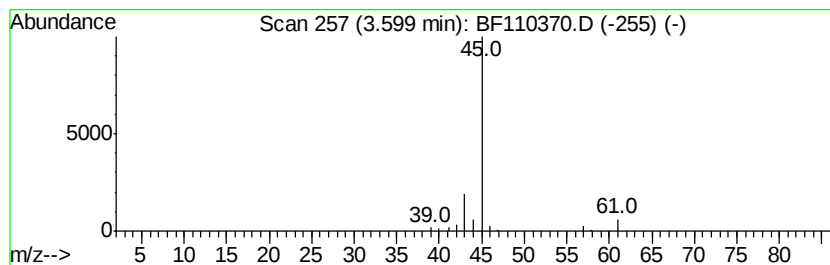
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TIC Integration Parameters: LSCINT.P

Peak Number 2 (S)-(+)-1,2-Propanediol Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.60	3.22 ng	90916	1,4-Dichlorobenzene-d4	6.96

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	(S)-(+)-1,2-Propanediol	76	C3H8O2	004254-15-3	74
2		R-(-)-1,2-propanediol	76	C3H8O2	004254-14-2	64
3		2-Butanol	74	C4H10O	000078-92-2	9
4		Propylene Glycol	76	C3H8O2	000057-55-6	9
5		Isopropyl Alcohol	60	C3H8O	000067-63-0	9



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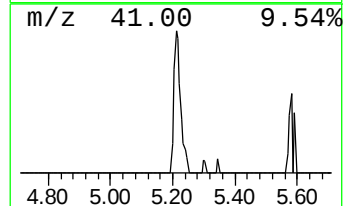
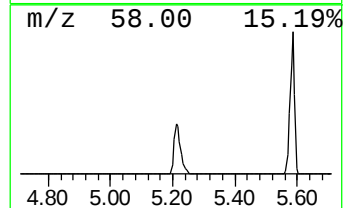
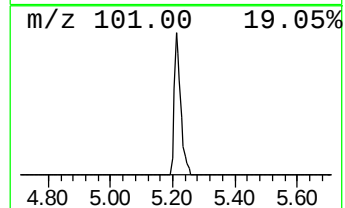
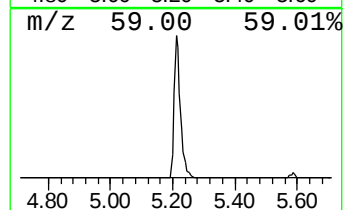
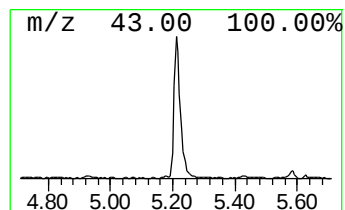
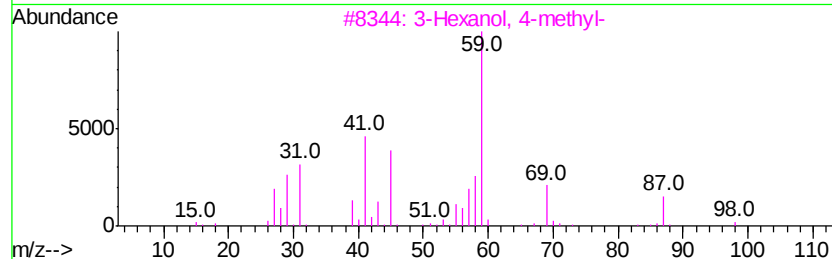
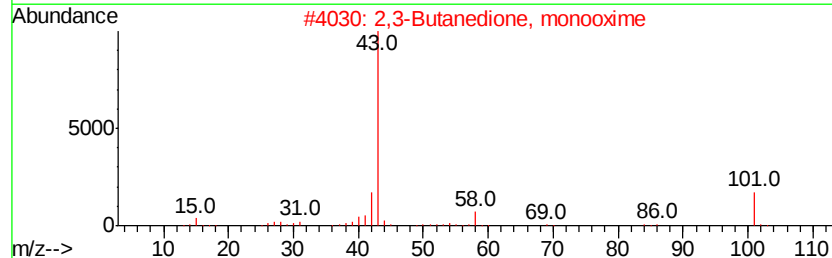
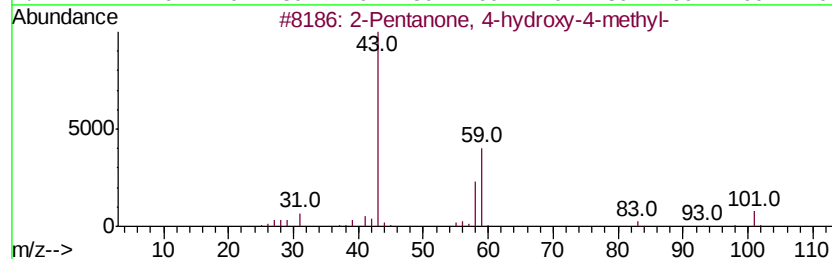
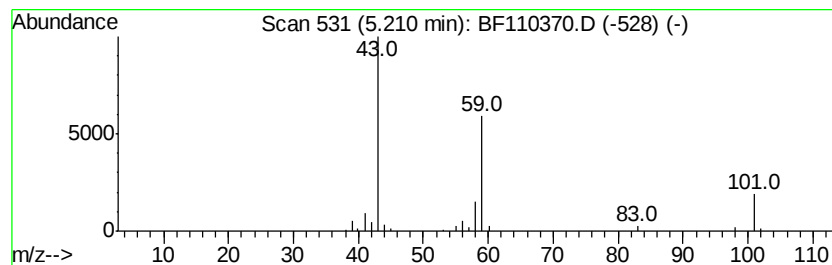
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.21	4.87 ng	137370	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	42
2		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	35
3		3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	28
4		Acetic acid, cyano-, 1,1-dimethyl-	141	C7H11NO2	001116-98-9	17
5		Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9



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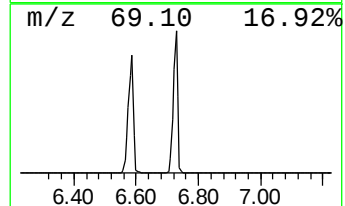
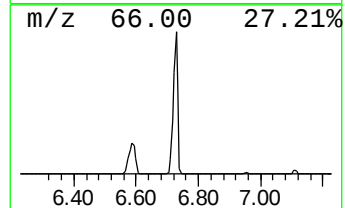
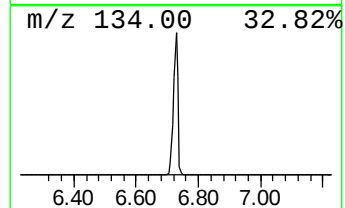
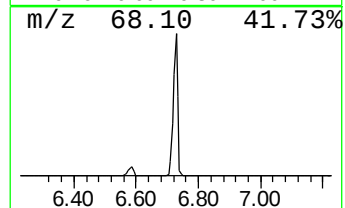
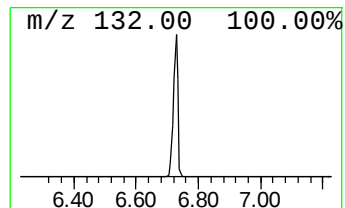
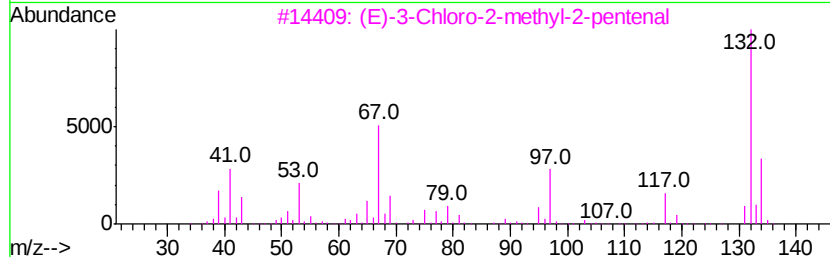
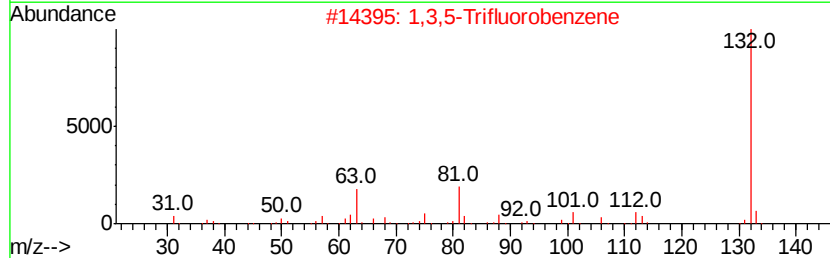
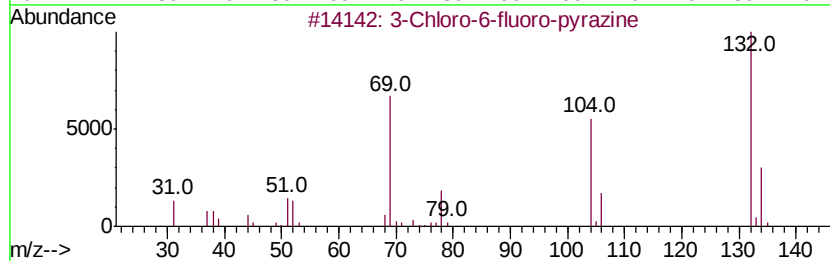
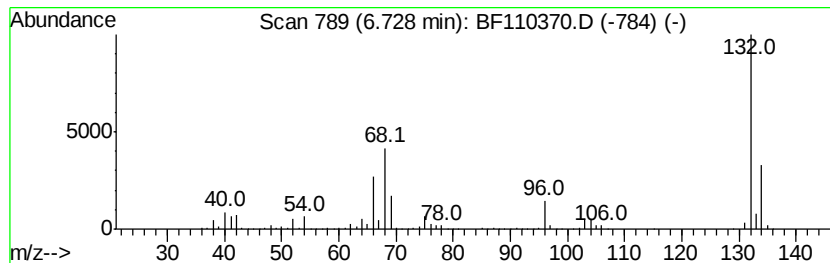
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 unknown6.73 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.73	106.28 ng	2997570	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	27
2		1,3,5-Trifluorobenzene	132	C6H3F3	000372-38-3	9
3		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	9
4		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
5		Benzene, 1-ethenyl-3,5-dimethyl-	132	C10H12	005379-20-4	9



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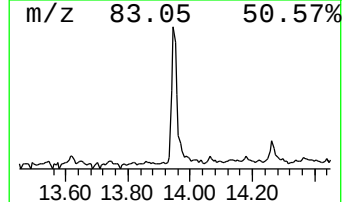
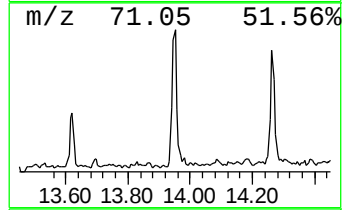
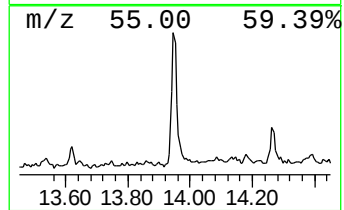
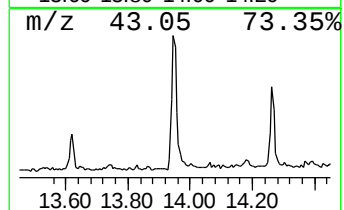
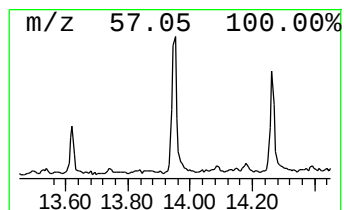
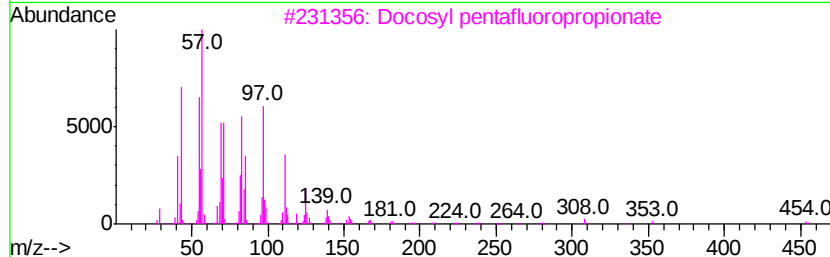
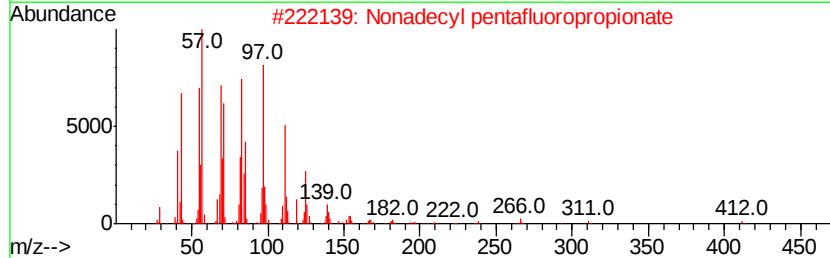
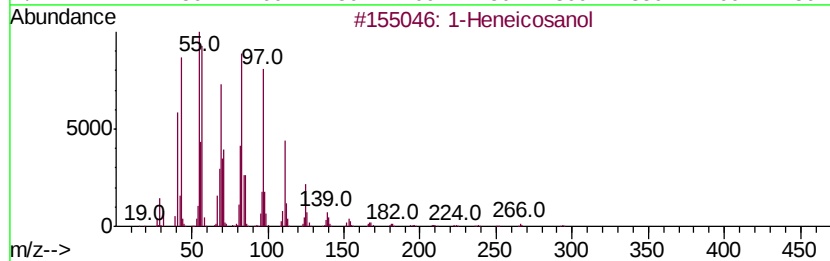
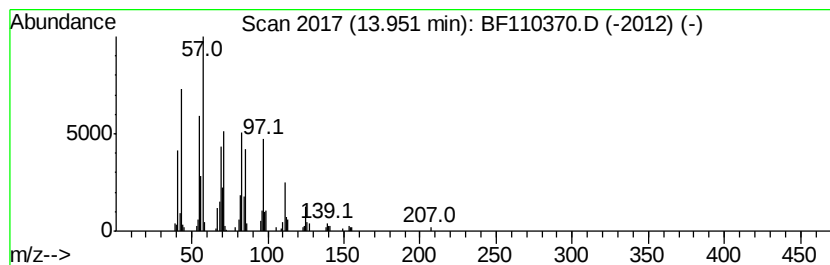
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 1-Heneicosanol Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD		R.T.		
13.95	9.03 ng	220774	Chrysene-d12		14.13		
Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Heneicosanol	312	C21H44O	015594-90-8	93
2			Nonadecyl pentafluoropropionate	430	C22H39F5O2	1000351-88-8	91
3			Docosyl pentafluoropropionate	472	C25H45F5O2	1000351-80-9	91
4			Octadecyl trifluoroacetate	366	C20H37F3O2	079392-43-1	91
5			Heneicosyl heptafluorobutyrate	508	C25H43F7O2	1000351-83-8	91



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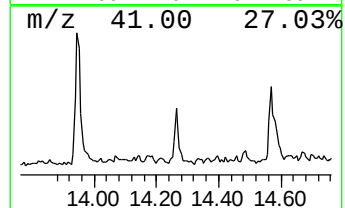
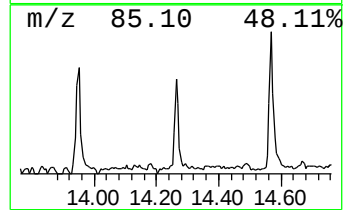
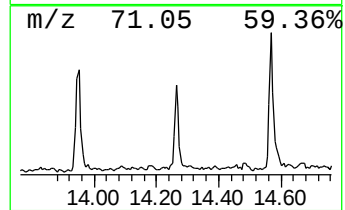
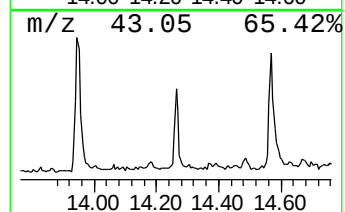
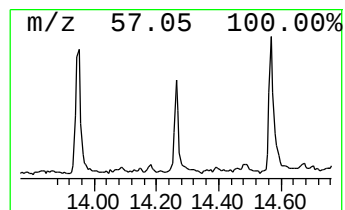
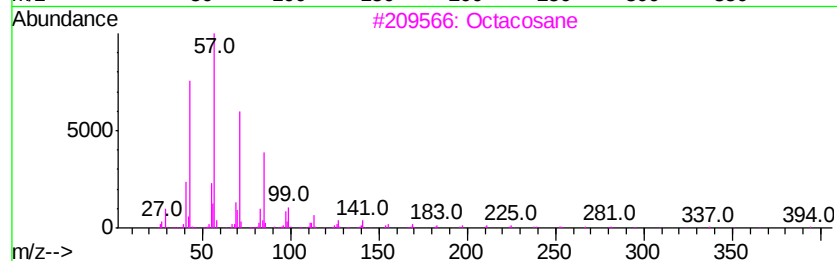
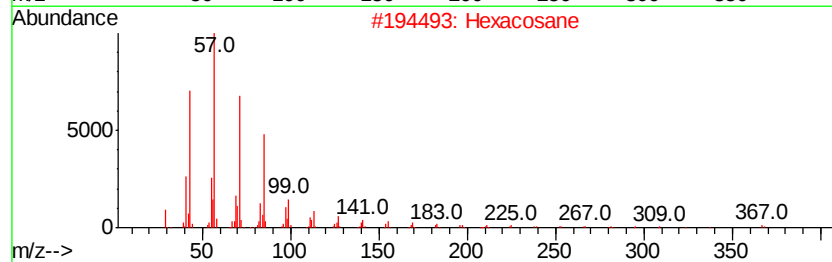
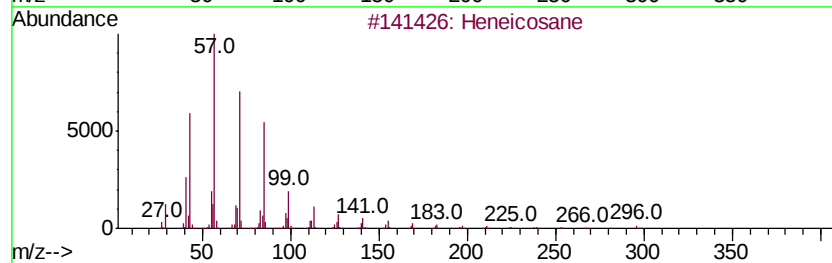
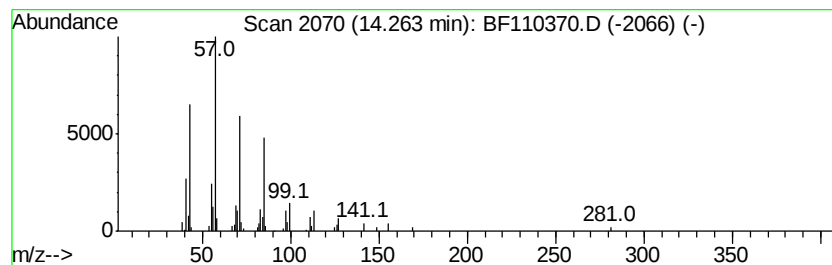
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Heneicosane Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.26	2.84 ng	69420	Chrysene-d12	14.13

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane		296	C21H44	000629-94-7	91
2	Hexacosane		366	C26H54	000630-01-3	91
3	Octacosane		394	C28H58	000630-02-4	91
4	Heptadecane		240	C17H36	000629-78-7	91
5	2-methyloctacosane		408	C29H60	1000376-72-8	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF110118\
Data File : BF110370.D
Acq On : 1 Nov 2018 15:55
Operator : JU/SJ
Sample : J5741-55
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TP-14

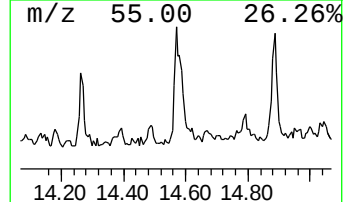
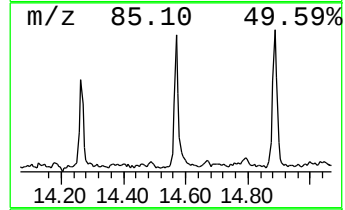
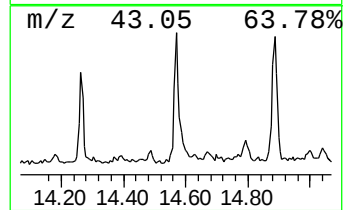
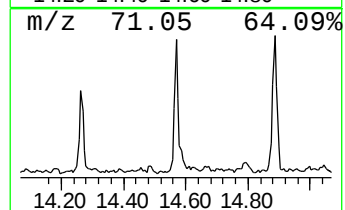
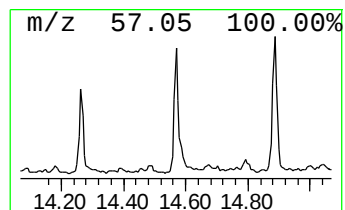
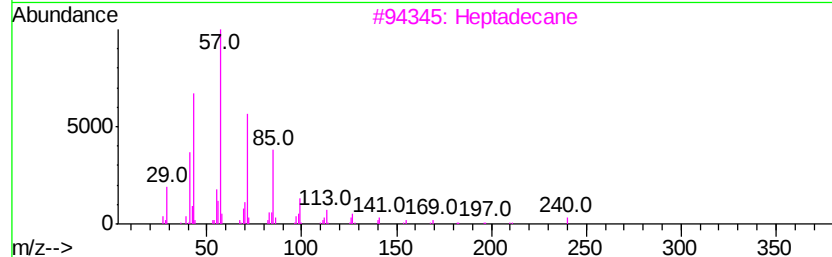
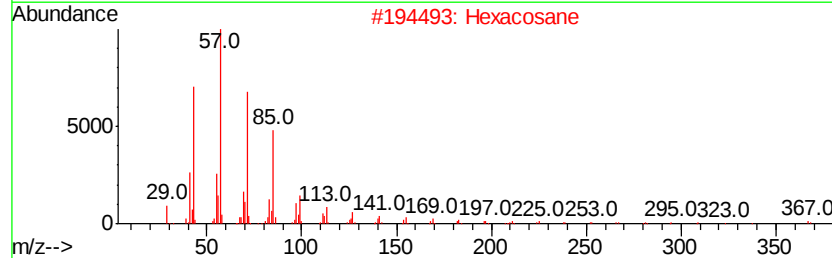
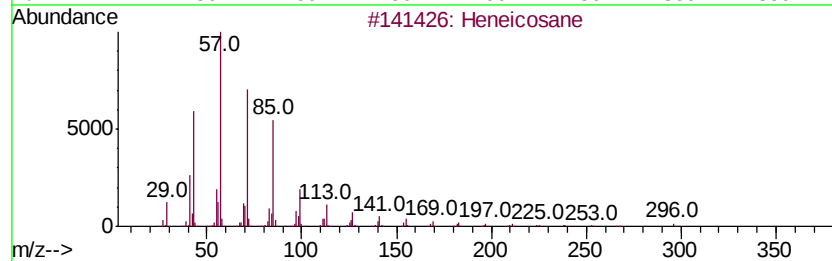
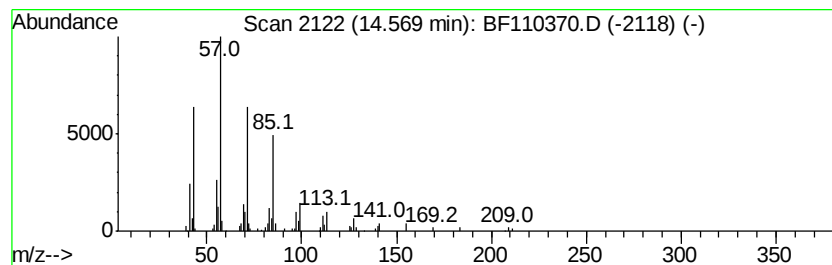
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Heptadecane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.57	6.13 ng	150012	Chrysene-d12	14.13

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heneicosane	296	C21H44	000629-94-7	91
2		Hexacosane	366	C26H54	000630-01-3	91
3		Heptadecane	240	C17H36	000629-78-7	91
4		Tetracosane	338	C24H50	000646-31-1	90
5		Hexadecane	226	C16H34	000544-76-3	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF110118\
 Data File : BF110370.D
 Acq On : 1 Nov 2018 15:55
 Operator : JU/SJ
 Sample : J5741-55
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

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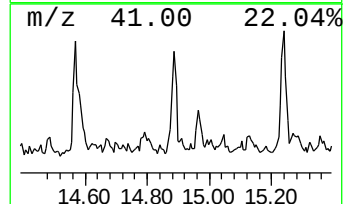
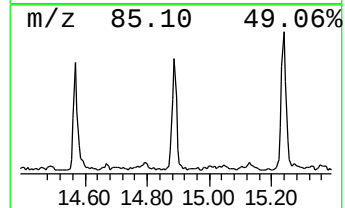
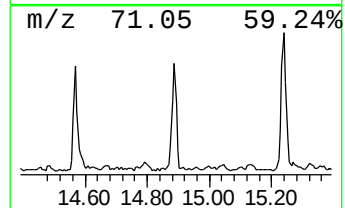
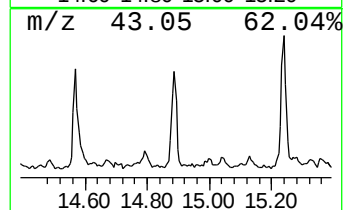
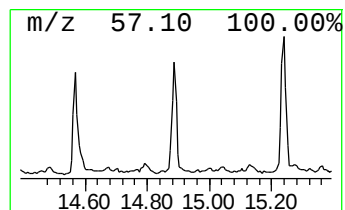
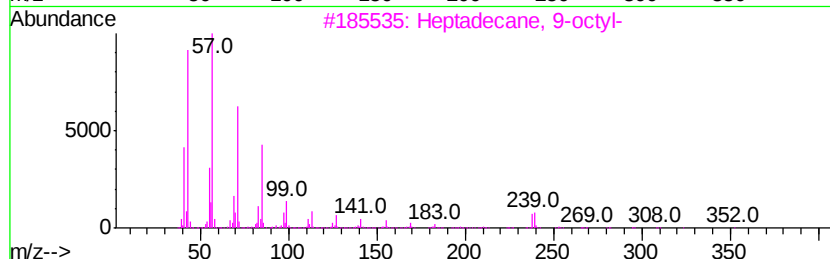
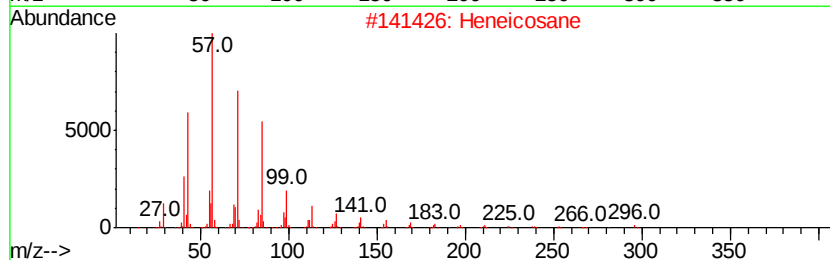
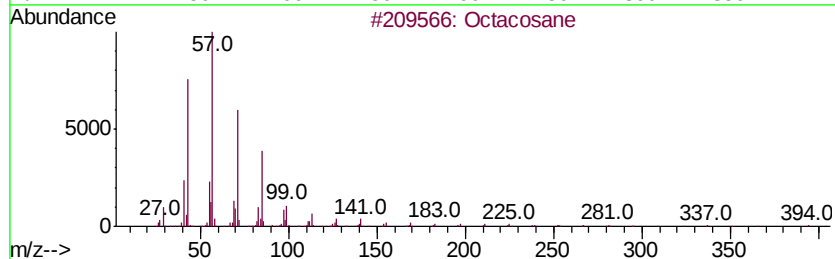
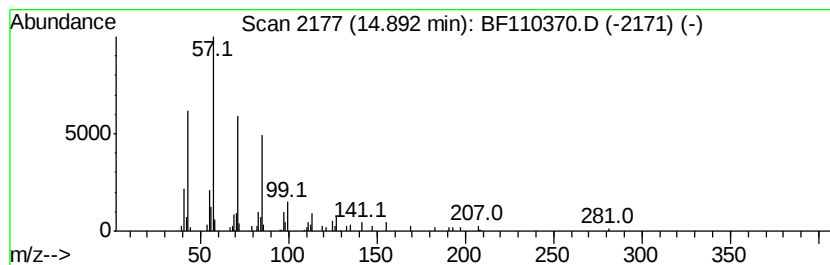
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

 Peak Number 9 Octacosane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.89	4.96 ng	121343	Chrysene-d12	14.13

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	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	Hentriacontane		437	C31H64	000630-04-6	91
5	Hexacosane		366	C26H54	000630-01-3	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF110118\
Data File : BF110370.D
Acq On : 1 Nov 2018 15:55
Operator : JU/SJ
Sample : J5741-55
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TP-14

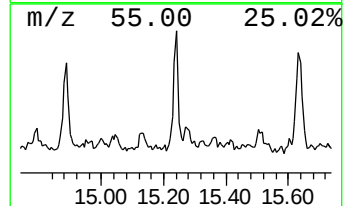
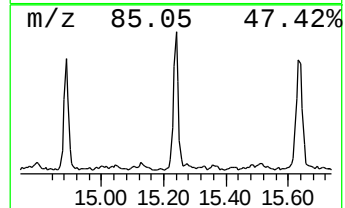
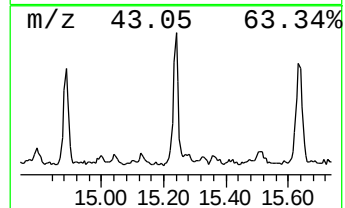
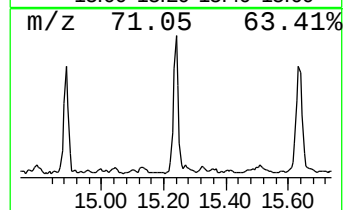
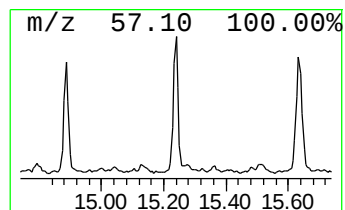
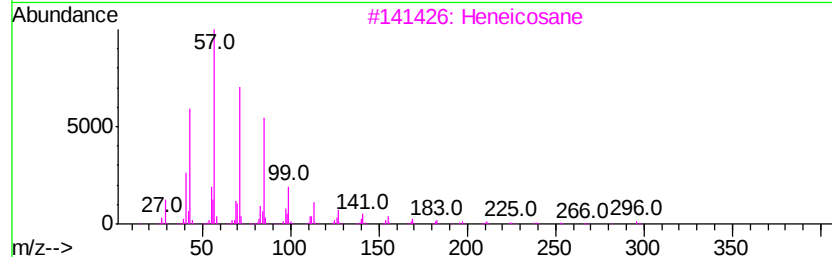
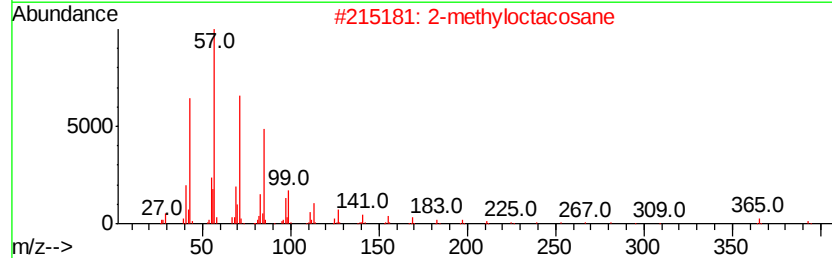
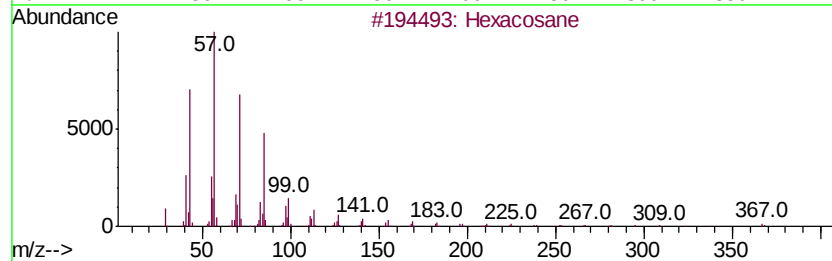
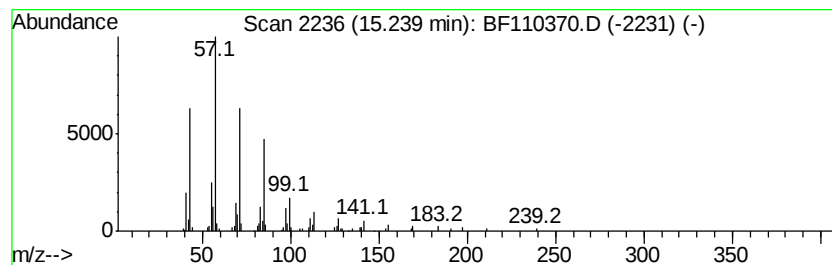
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Hexacosane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.24	4.84 ng	158696	Perylene-d12	15.66

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	91
3		Heneicosane	296	C21H44	000629-94-7	91
4		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
5		Triacontane	422	C30H62	000638-68-6	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF110118\
Data File : BF110370.D
Acq On : 1 Nov 2018 15:55
Operator : JU/SJ
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ClientSampleId :
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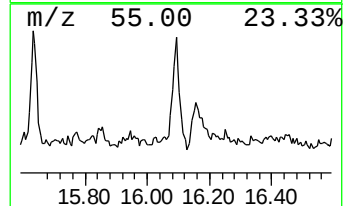
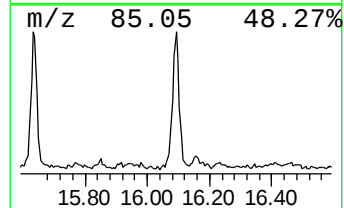
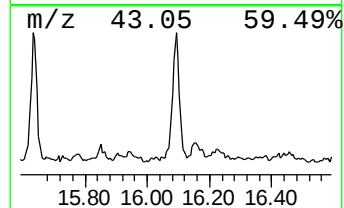
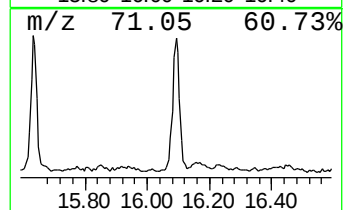
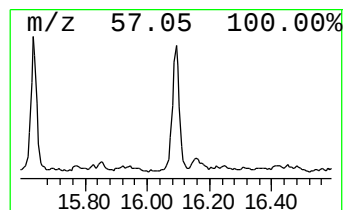
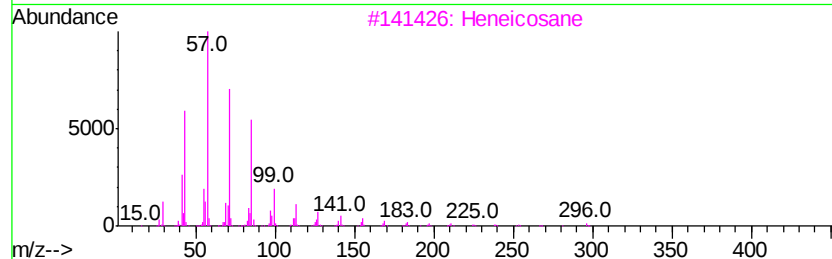
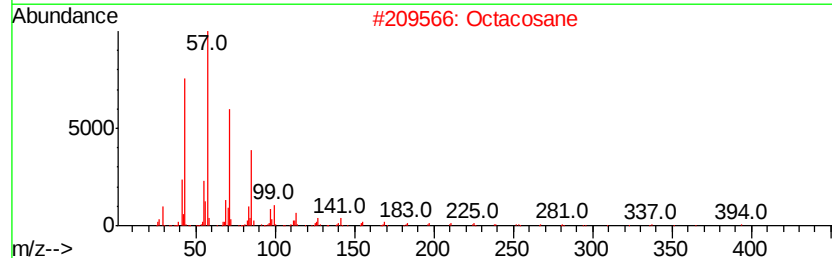
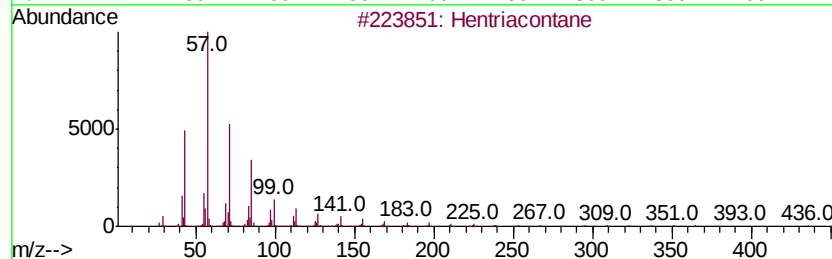
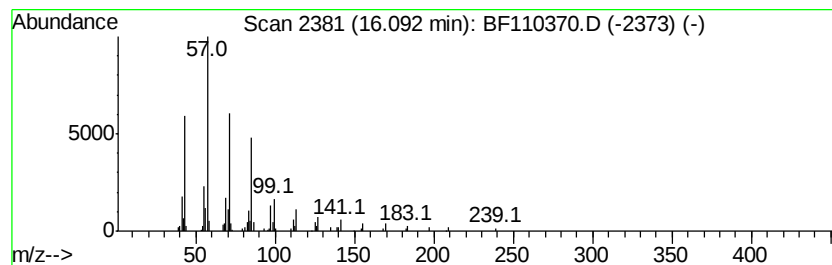
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Hentriacontane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.09	5.57 ng	182594	Perylene-d12	15.66

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hentriacontane	437	C31H64	000630-04-6	91
2			Octacosane	394	C28H58	000630-02-4	91
3			Heneicosane	296	C21H44	000629-94-7	91
4			Docosane	310	C22H46	000629-97-0	91
5			Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF110118\
Data File : BF110370.D
Acq On : 1 Nov 2018 15:55
Operator : JU/SJ
Sample : J5741-55
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TP-14

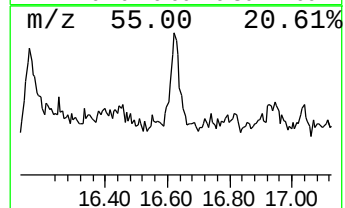
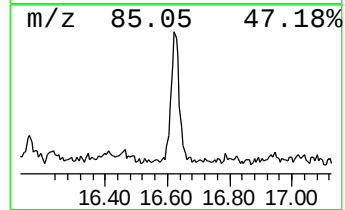
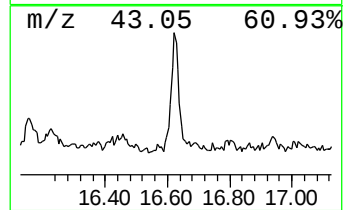
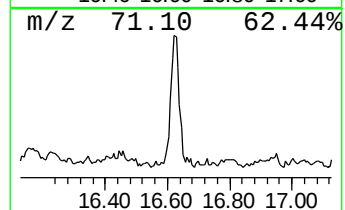
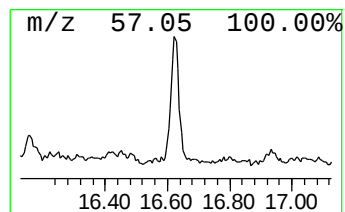
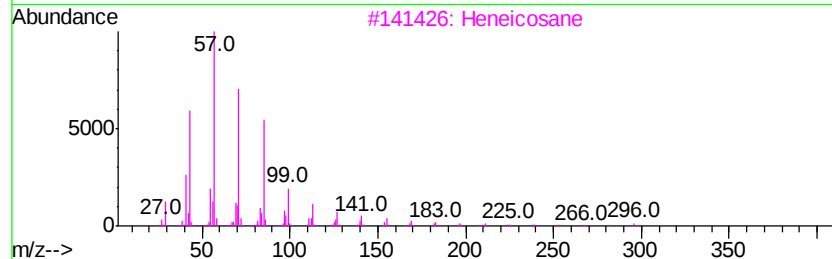
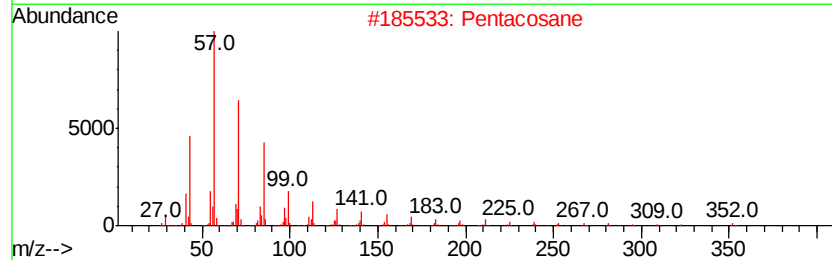
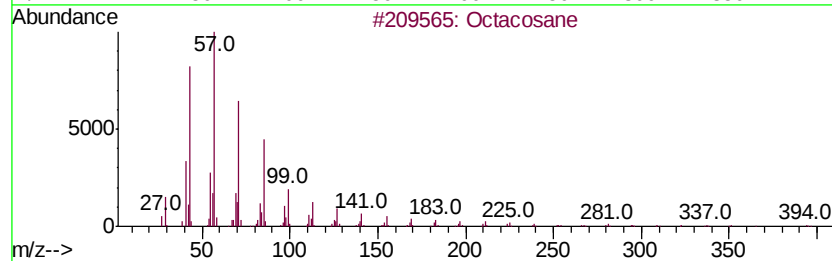
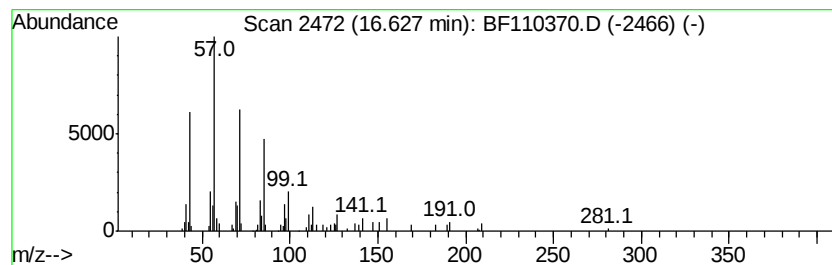
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Pentacosane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.63	3.03 ng	99507	Perylene-d12	15.66

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octacosane	394	C28H58	000630-02-4	90
2			Pentacosane	352	C25H52	000629-99-2	90
3			Heneicosane	296	C21H44	000629-94-7	87
4			Heptacosane	380	C27H56	000593-49-7	87
5			Hexacosane	366	C26H54	000630-01-3	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF110118\
Data File : BF110370.D
Acq On : 1 Nov 2018 15:55
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ALS Vial : 11 Sample Multiplier: 1

Instrument :
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ClientSampled :
TP-14

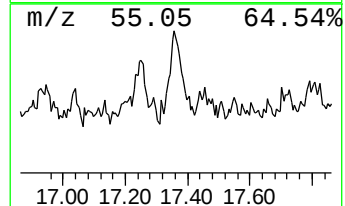
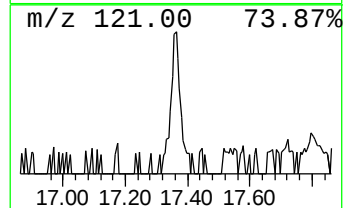
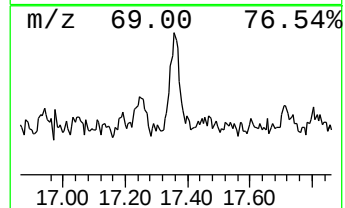
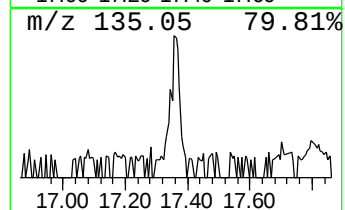
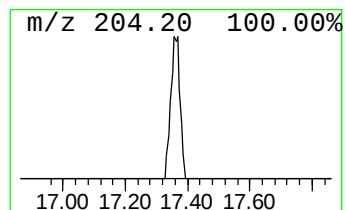
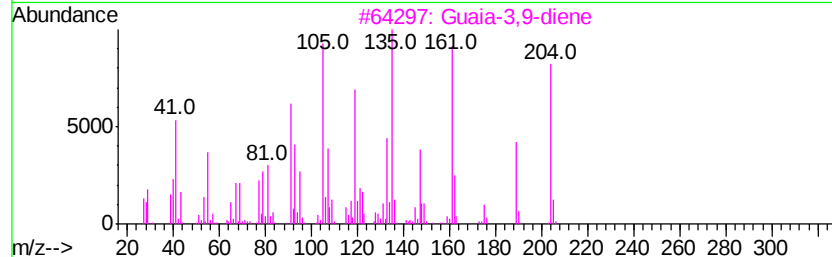
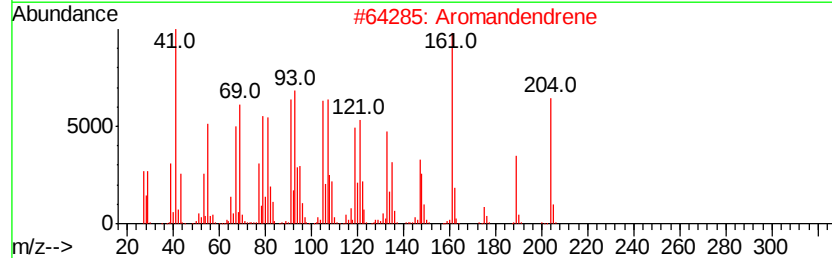
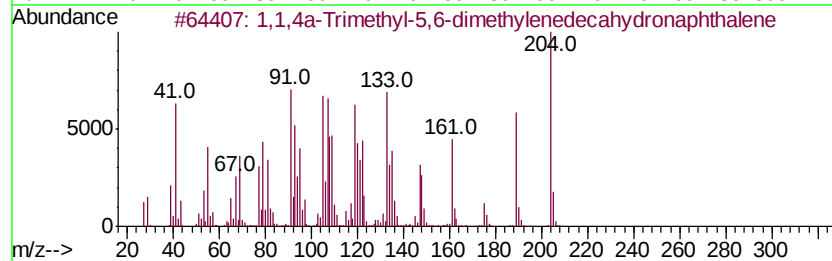
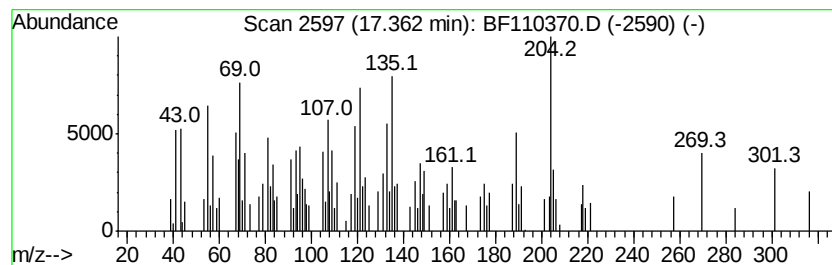
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 1,1,4a-Trimethyl-5,6-dimeth... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.36	4.32 ng	141733	Perylene-d12	15.66

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,1,4a-Trimethyl-5,6-dimethylene...	204	C15H24	1000193-60-8	78
2			Aromandendrene	204	C15H24	000489-39-4	64
3			Guaia-3,9-diene	204	C15H24	000489-83-8	62
4			2H-Cyclopentacyclooctene, 4,5,6,...	204	C15H24	1000221-85-8	55
5			1H-3a,7-Methanoazulene, octahydr...	204	C15H24	000508-55-4	53



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF110118\
 Data File : BF110370.D
 Acq On : 1 Nov 2018 15:55
 Operator : JU/SJ
 Sample : J5741-55
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF102318.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
Butane, 2-methoxy...	2.32	21.6	ng	608053	1	6.96	564116	20.0
(S)-(+)-1,2-Propa...	3.60	3.2	ng	90916	1	6.96	564116	20.0
2-Pentanone, 4-hy...	5.21	4.9	ng	137370	1	6.96	564116	20.0
unknown6.73	6.73	106.3	ng	2997570	1	6.96	564116	20.0
1-Heneicosanol	13.95	9.0	ng	220774	5	14.13	489053	20.0
Heneicosane	14.26	2.8	ng	69420	5	14.13	489053	20.0
Heptadecane	14.57	6.1	ng	150012	5	14.13	489053	20.0
Octacosane	14.89	5.0	ng	121343	5	14.13	489053	20.0
Hexacosane	15.24	4.8	ng	158696	6	15.66	655935	20.0
Hentriacontane	16.09	5.6	ng	182594	6	15.66	655935	20.0
Pentacosane	16.63	3.0	ng	99507	6	15.66	655935	20.0
1,1,4a-Trimethyl-...	17.36	4.3	ng	141733	6	15.66	655935	20.0