

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF022819\
 Data File : BF112781.D
 Acq On : 1 Mar 2019 00:27
 Operator : JU/SJ
 Sample : K1651-04
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CON-327

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-BF022719.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	2.987	150	153	163	rBV2	36143	63753	1.10%	0.229%
2	3.269	194	201	210	rBV	69684	135260	2.34%	0.486%
3	4.275	364	372	376	rBV	3710032	5624531	97.45%	20.218%
4	4.940	476	485	488	rBV	3523432	5771598	100.00%	20.747%
5	5.340	549	553	557	rBV	59960	61903	1.07%	0.223%
6	6.363	723	727	738	rBV	661023	551137	9.55%	1.981%
7	6.504	748	751	758	rVB	138144	115506	2.00%	0.415%
8	6.740	788	791	795	rBV	684217	544920	9.44%	1.959%
9	6.898	814	818	821	rBV	1941716	1537255	26.63%	5.526%
10	7.057	842	845	849	rBV	107609	83154	1.44%	0.299%
11	7.298	882	886	889	rBV	1427393	1191539	20.64%	4.283%
12	8.022	1006	1009	1018	rBV	828174	715078	12.39%	2.570%
13	9.098	1188	1192	1202	rBV	2802044	2339405	40.53%	8.409%
14	9.222	1210	1213	1225	rVB	194416	167418	2.90%	0.602%
15	9.486	1256	1258	1265	rBV	127229	117741	2.04%	0.423%
16	9.769	1303	1306	1314	rBV	855281	796213	13.80%	2.862%
17	11.245	1554	1557	1560	rBV	914732	801358	13.88%	2.881%
18	11.322	1567	1570	1573	rVB	85888	72341	1.25%	0.260%
19	12.298	1733	1736	1741	rBV	254853	240648	4.17%	0.865%
20	12.451	1759	1762	1766	rBV	292310	256865	4.45%	0.923%
21	12.680	1796	1801	1803	rBV	260278	274091	4.75%	0.985%
22	12.827	1822	1826	1830	rVB	2474887	2148512	37.23%	7.723%
23	13.074	1865	1868	1871	rVB	71299	66057	1.14%	0.237%
24	13.733	1976	1980	1984	rBV2	299369	366958	6.36%	1.319%
25	13.869	1996	2003	2006	rBV2	693328	783493	13.57%	2.816%
26	13.898	2006	2008	2011	rVB	200184	156715	2.72%	0.563%
27	14.057	2032	2035	2038	rBV	135022	128623	2.23%	0.462%
28	14.363	2084	2087	2092	rBV	260280	224435	3.89%	0.807%
29	14.657	2134	2137	2141	rVB	234542	234615	4.06%	0.843%
30	14.880	2171	2175	2178	rBV	76211	106545	1.85%	0.383%
31	14.980	2188	2192	2202	rVB	320585	423139	7.33%	1.521%
32	15.280	2238	2243	2248	rVV	515970	639040	11.07%	2.297%
33	15.333	2248	2252	2261	rVB	205896	274232	4.75%	0.986%
34	15.521	2280	2284	2289	rVB	119365	158742	2.75%	0.571%

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Integration Parameters: rteint.p
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

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Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	15.745	2317	2322	2325	rBV	195284	282392	4.89%	1.015%
36	15.786	2325	2329	2340	rVB2	113322	204920	3.55%	0.737%
37	16.215	2398	2402	2413	rVB	88852	159130	2.76%	0.572%

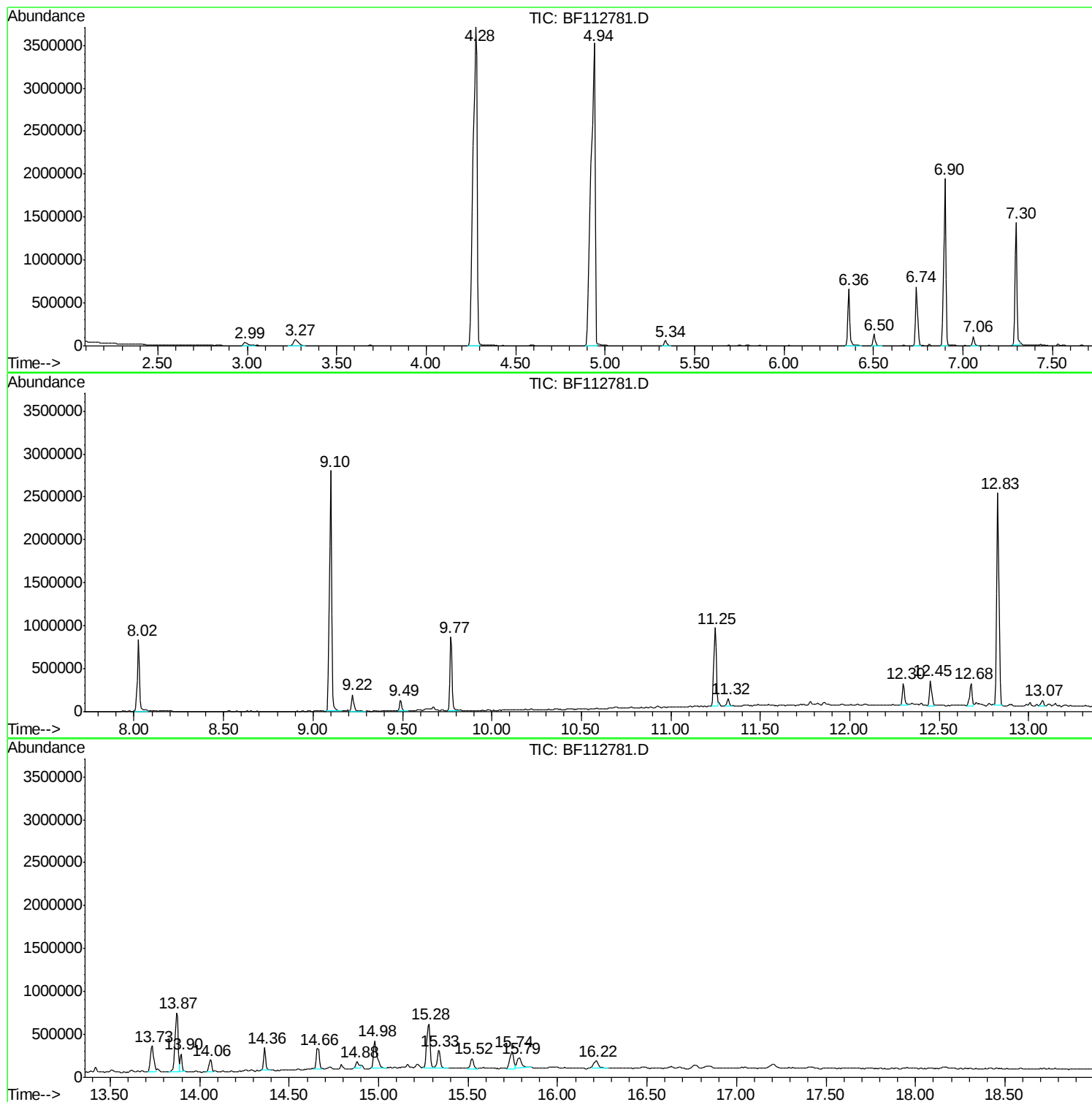
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF022719.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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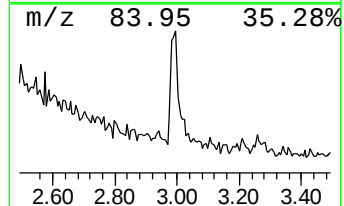
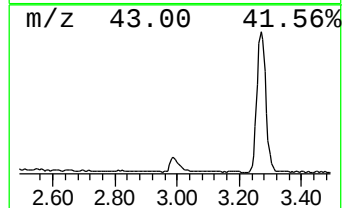
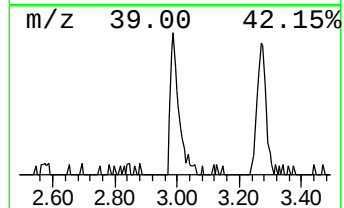
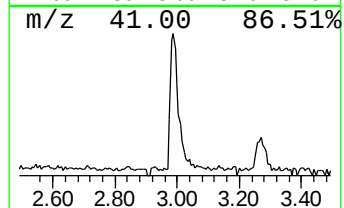
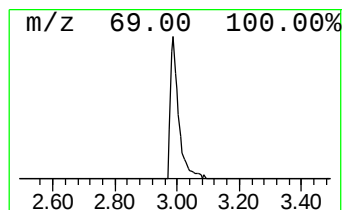
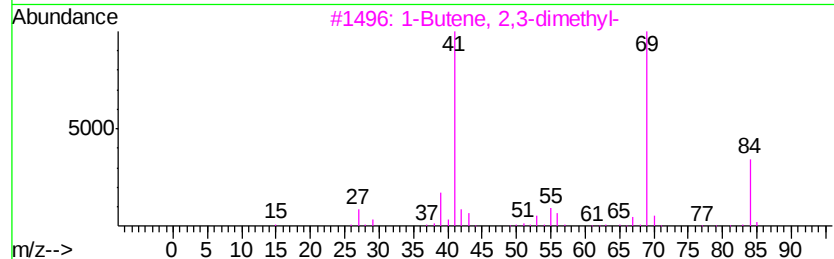
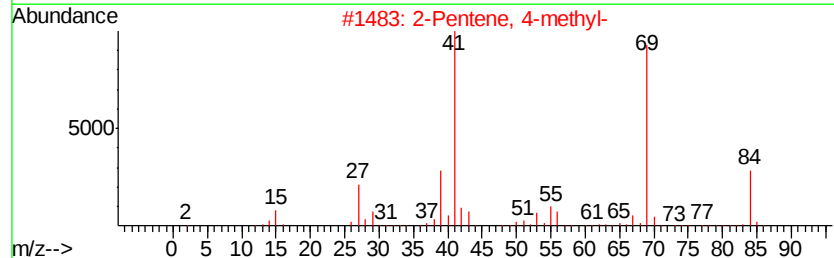
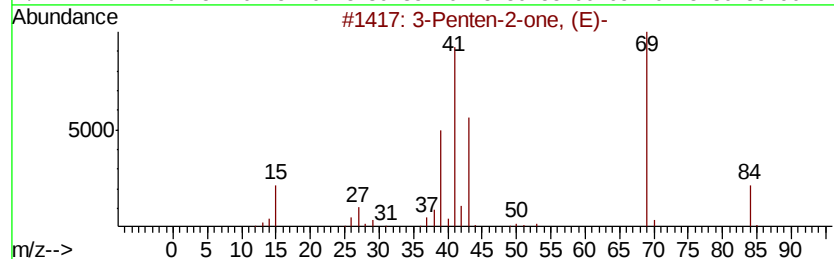
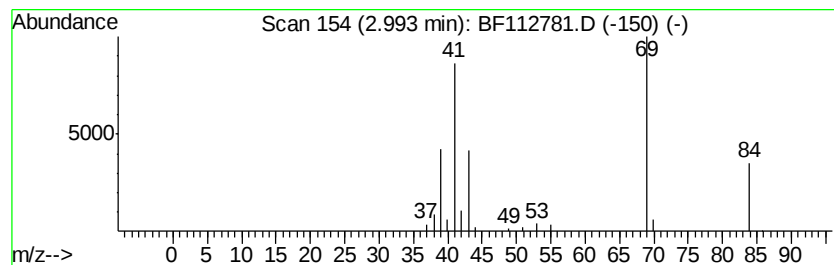
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF022719.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 3-Penten-2-one, (E)- Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.99	2.34 ng	63753	1,4-Dichlorobenzene-d4	6.74

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Penten-2-one, (E)-	84	C5H8O	003102-33-8	83
2			2-Pentene, 4-methyl-	84	C6H12	004461-48-7	53
3			1-Butene, 2,3-dimethyl-	84	C6H12	000563-78-0	53
4			2-Pentene, 4-methyl-, (E)-	84	C6H12	000674-76-0	52
5			3-Penten-2-one	84	C5H8O	000625-33-2	50



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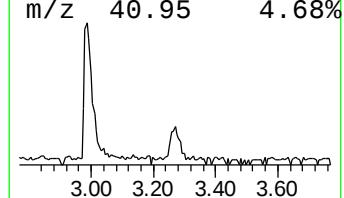
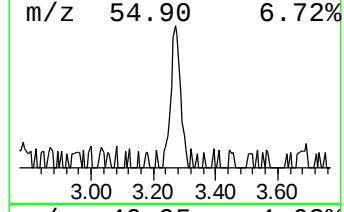
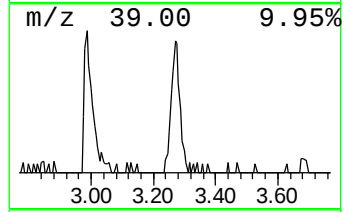
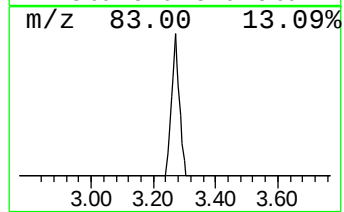
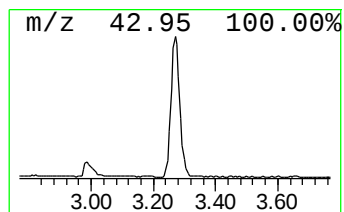
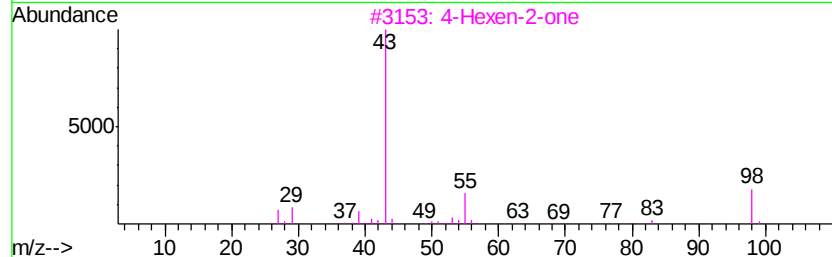
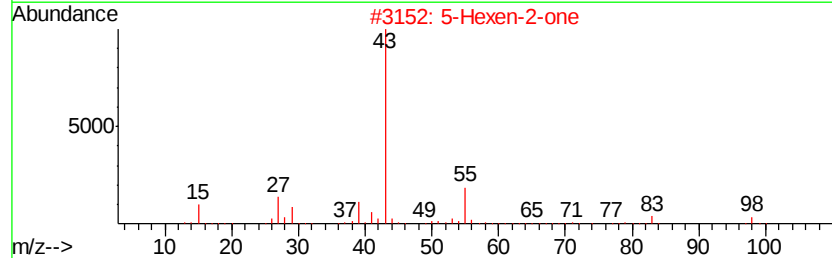
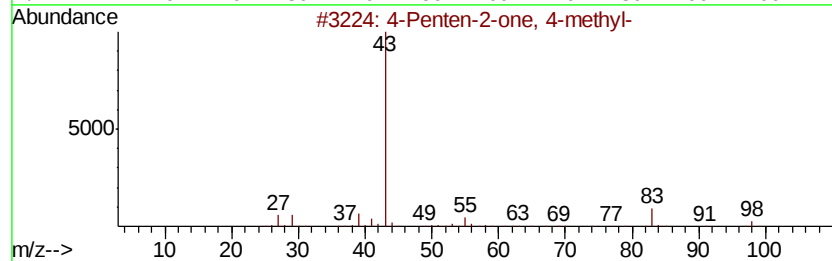
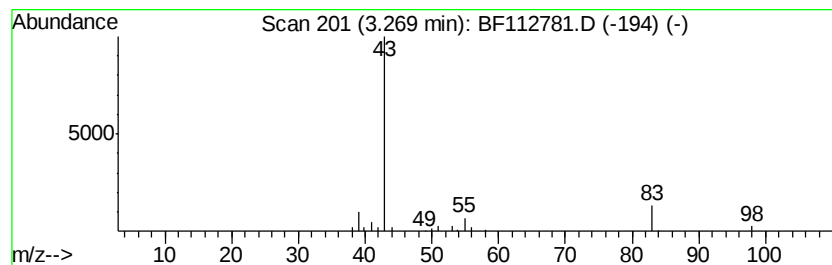
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 4-Penten-2-one, 4-methyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.27	4.96 ng	135260	1,4-Dichlorobenzene-d4	6.74

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4-Penten-2-one, 4-methyl-	98	C6H10O	003744-02-3	86
2			5-Hexen-2-one	98	C6H10O	000109-49-9	35
3			4-Hexen-2-one	98	C6H10O	025659-22-7	9
4			2-Pentanone, 3-methylene-	98	C6H10O	004359-77-7	9
5			4-Penten-2-one, 3-methyl-	98	C6H10O	000758-87-2	9



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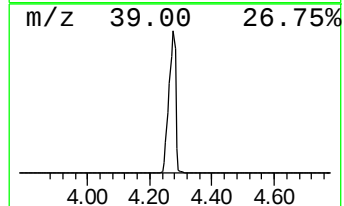
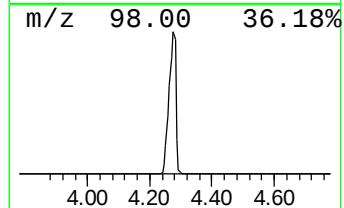
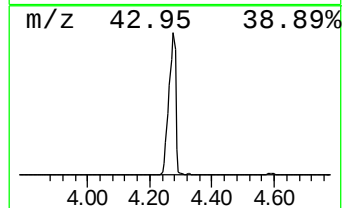
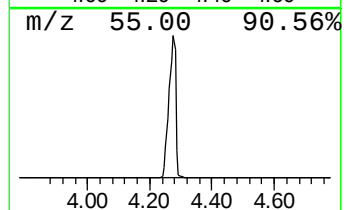
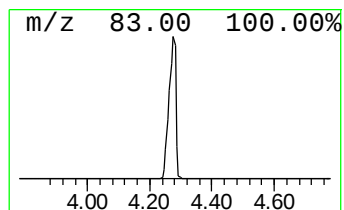
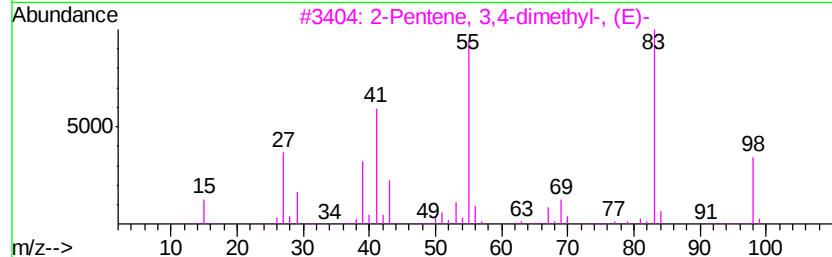
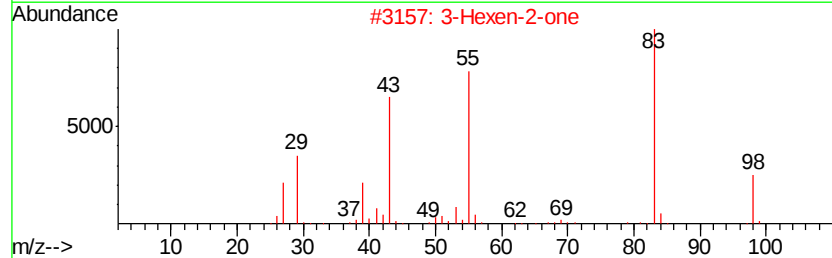
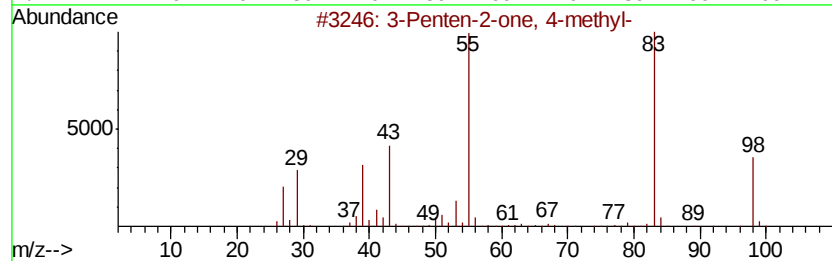
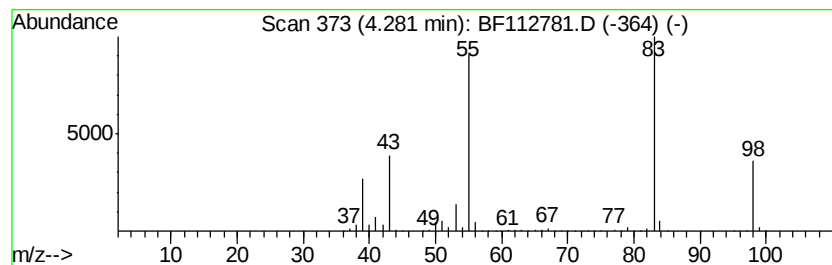
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.28	206.44 ng	5624530	1,4-Dichlorobenzene-d4	6.74

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Penten-2-one, 4-methyl-	98	C6H10O	000141-79-7	91
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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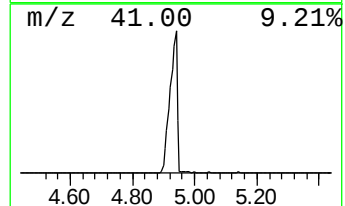
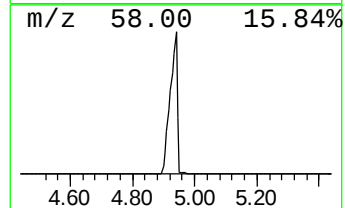
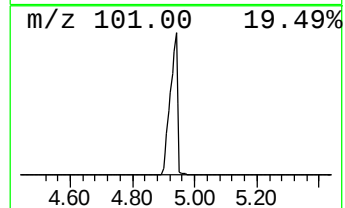
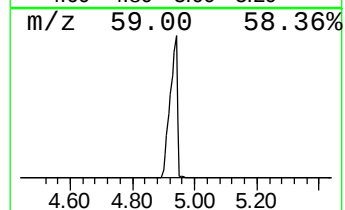
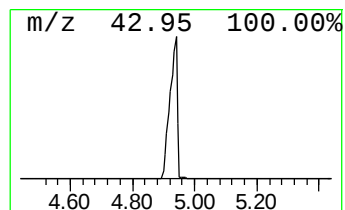
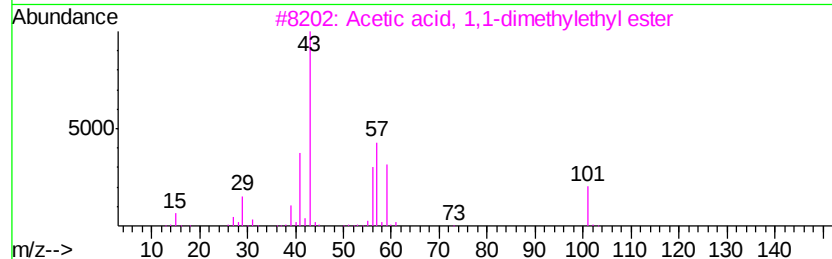
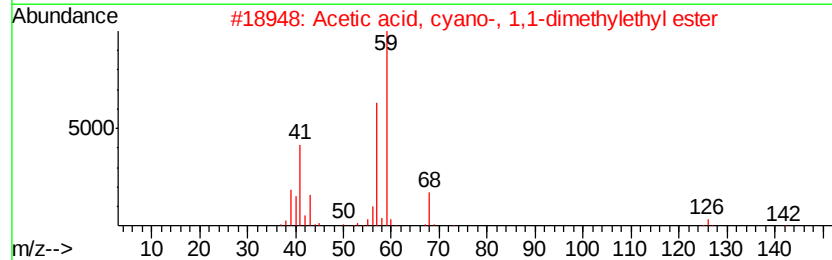
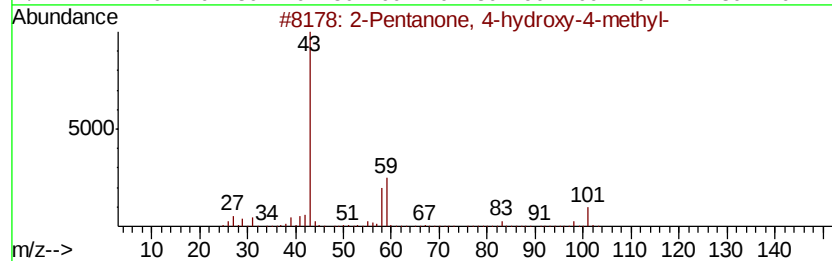
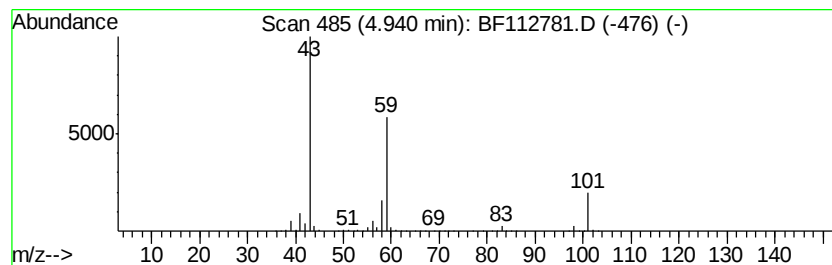
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.94	211.83 ng	5771600	1,4-Dichlorobenzene-d4	6.74

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2		Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	25
3		Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	9
4		Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	9
5		Propane, 1-ethoxy-2-methyl-	102	C6H14O	000627-02-1	9



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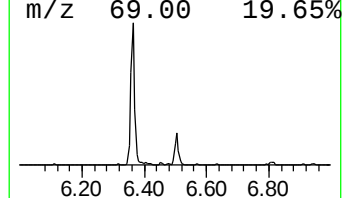
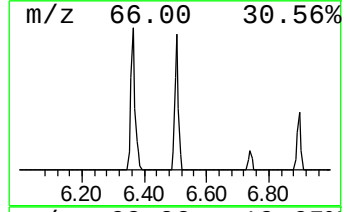
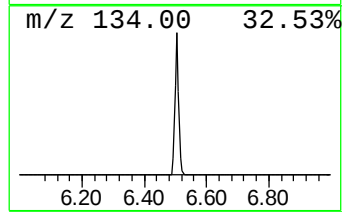
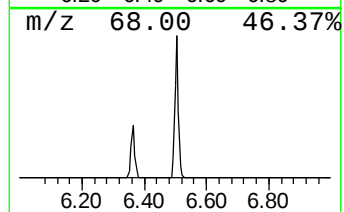
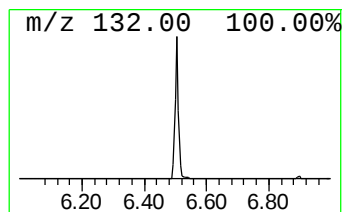
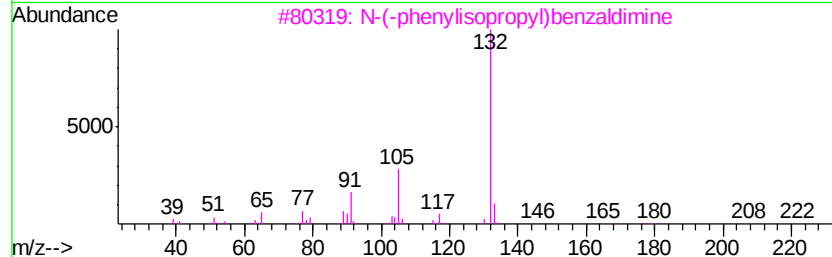
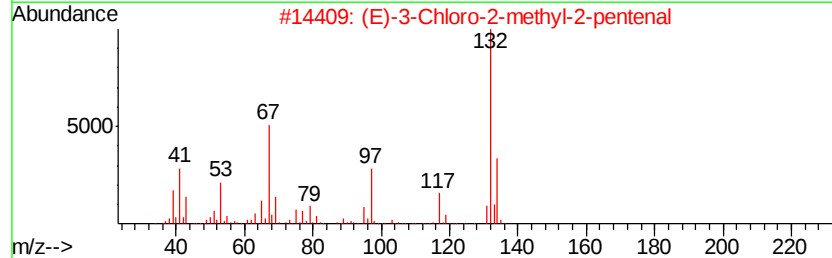
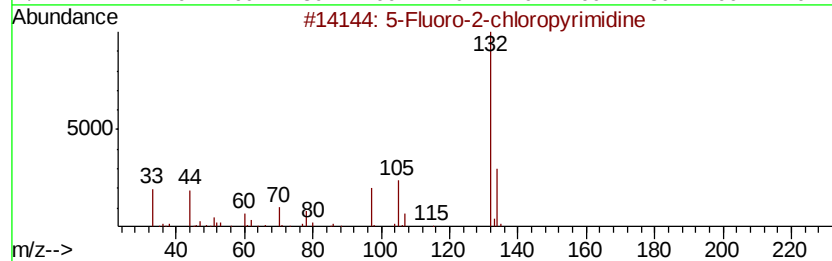
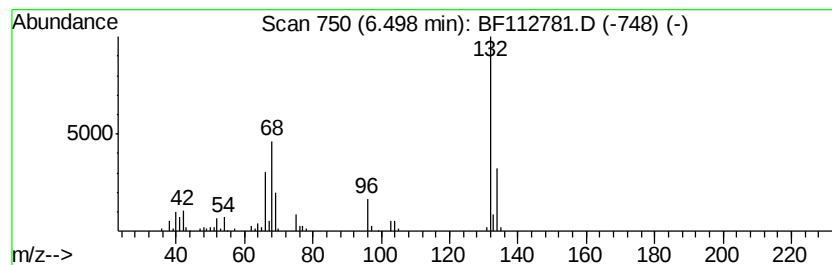
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF022719.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 unknown6.50 Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.50	4.24 ng	115506	1,4-Dichlorobenzene-d4	6.74

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	12
3		N-(-phenylisopropyl)benzalimine	223	C16H17N	1000379-01-3	10
4		5-Aminoindole	132	C8H8N2	005192-03-0	10
5		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF022819\
Data File : BF112781.D
Acq On : 1 Mar 2019 00:27
Operator : JU/SJ
Sample : K1651-04
Misc :
ALS Vial : 17 Sample Multiplier: 1

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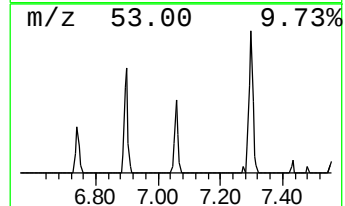
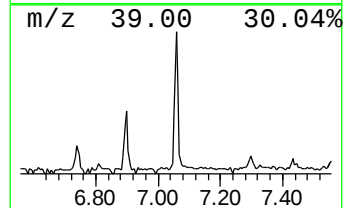
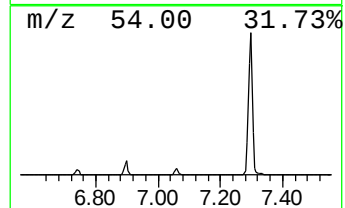
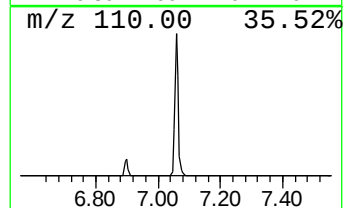
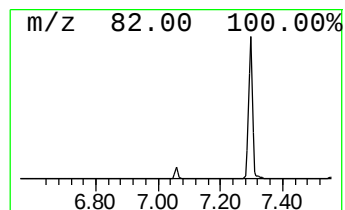
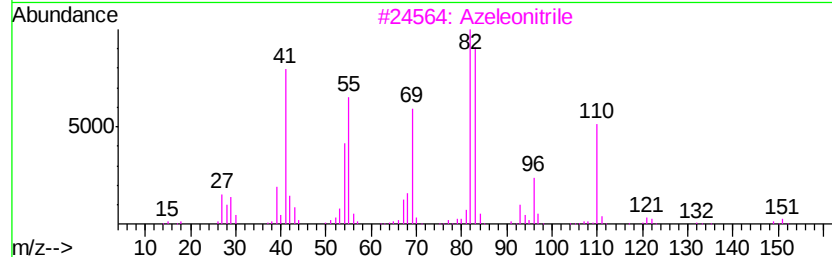
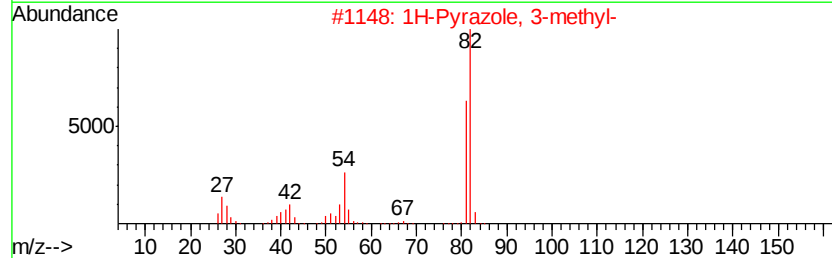
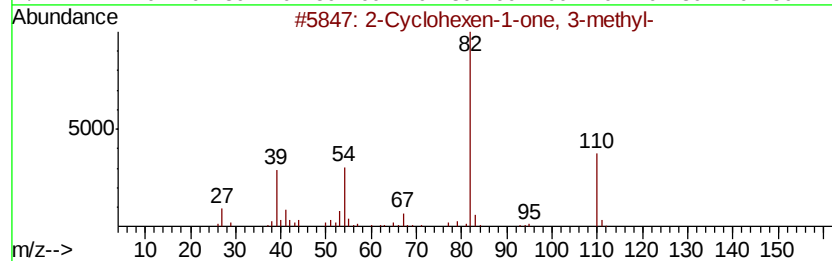
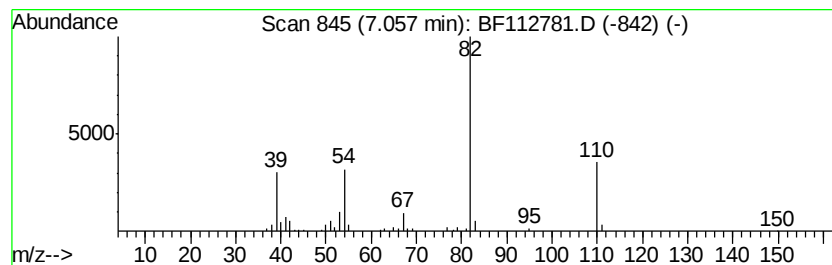
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 2-Cyclohexen-1-one, 3-methyl- Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.06	3.05 ng	83154	1,4-Dichlorobenzene-d4	6.74

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Cyclohexen-1-one, 3-methyl-	110	C7H10O	001193-18-6	91
2			1H-Pyrazole, 3-methyl-	82	C4H6N2	001453-58-3	59
3			Azeleoneitrile	150	C9H14N2	001675-69-0	56
4			1H-Imidazole, 4-methyl-	82	C4H6N2	000822-36-6	42
5			1'-Hydroxy-4,3'-dimethyl-bicyclo...	220	C14H20O2	1000189-12-5	39



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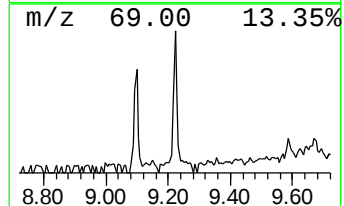
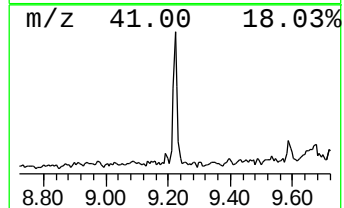
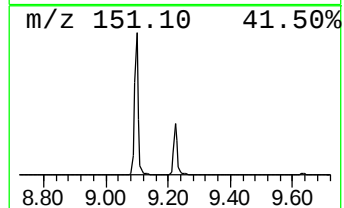
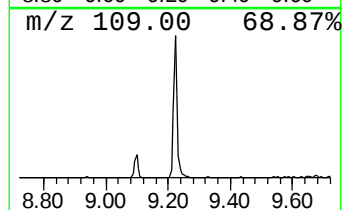
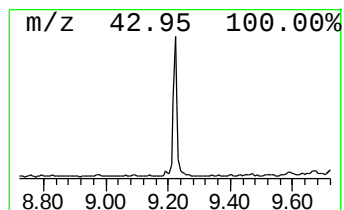
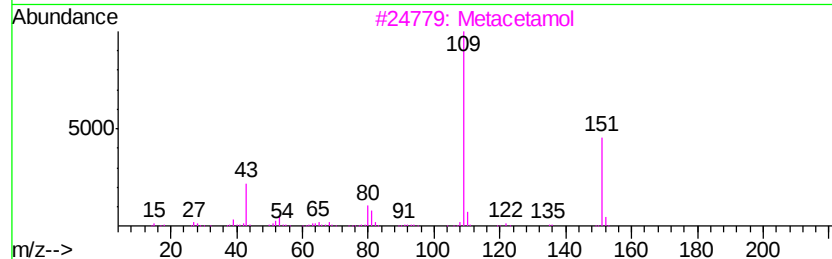
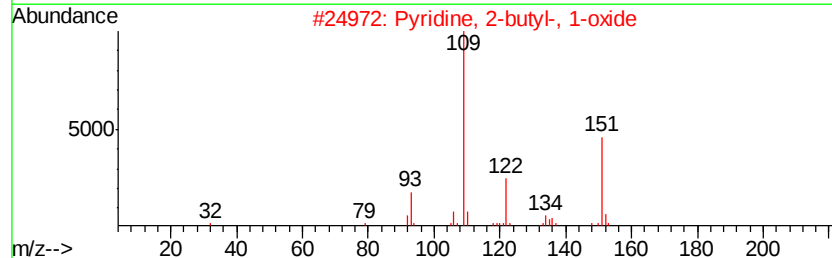
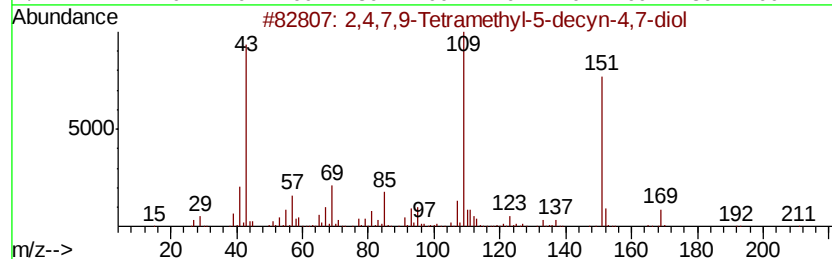
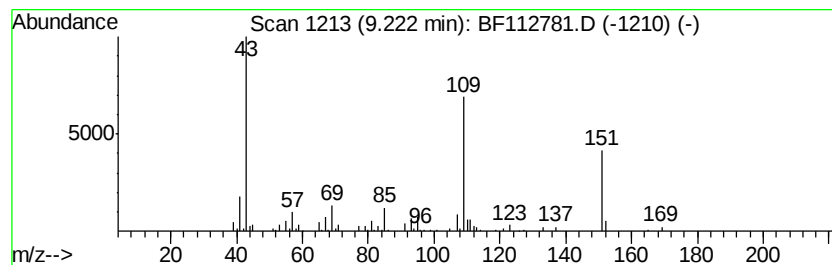
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 2,4,7,9-Tetramethyl-5-decyn... Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.22	4.21 ng	167418	Acenaphthene-d10	9.77

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2,4,7,9-Tetramethyl-5-decyn-4,7-...	226	C14H26O2	000126-86-3	90
2		Pyridine, 2-butyl-, 1-oxide	151	C9H13NO	031396-32-4	59
3		Metacetamol	151	C8H9NO2	000621-42-1	58
4		m-Isopropoxvaniline	151	C9H13NO	041406-00-2	42
5		Pyridine, 3-butyl-, 1-oxide	151	C9H13NO	031396-33-5	36



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF022819\
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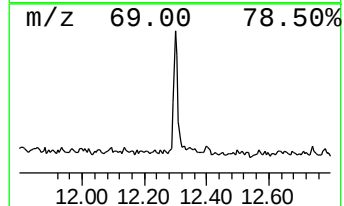
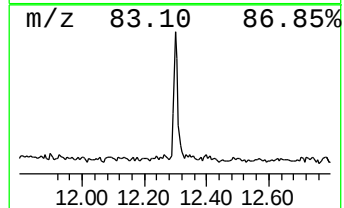
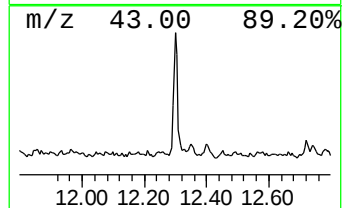
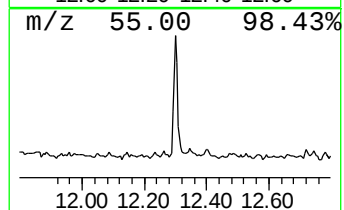
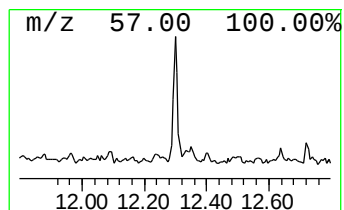
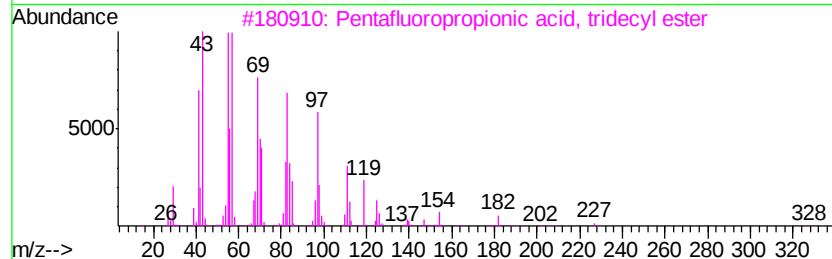
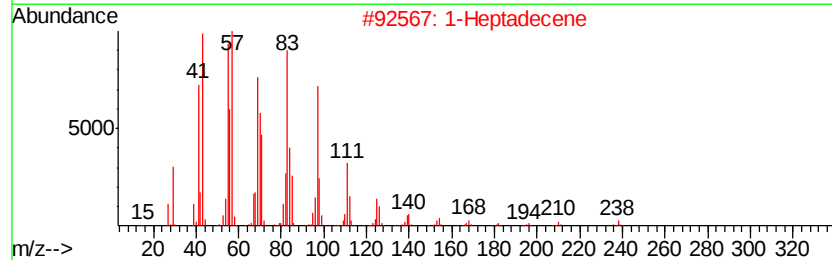
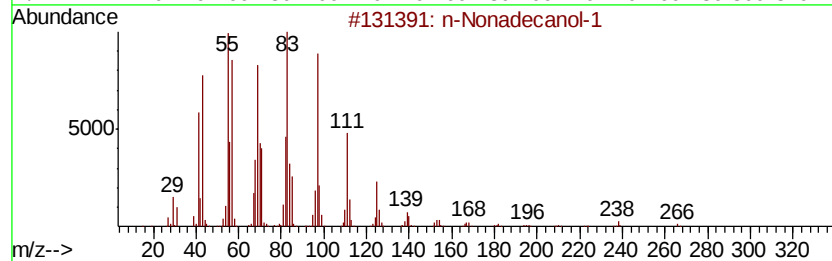
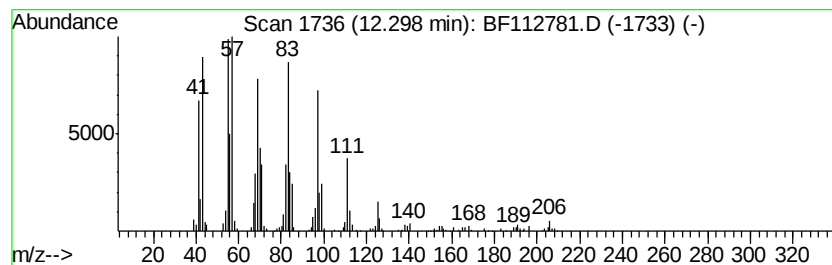
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 n-Nonadecanol-1 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.30	6.01 ng	240648	Phenanthrene-d10	11.25

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Nonadecanol-1	284	C19H40O	001454-84-8	91
2	1-Heptadecene	238	C17H34	006765-39-5	91
3	Pentafluoropropionic acid, tridecyl ester	346	C16H27F5O2	959261-22-4	91
4	4-Heptafluorobutyltridecylhexadecane	438	C20H33F7O2	1000282-97-2	91
5	1-Heneicosanol	312	C21H44O	015594-90-8	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF022819\
Data File : BF112781.D
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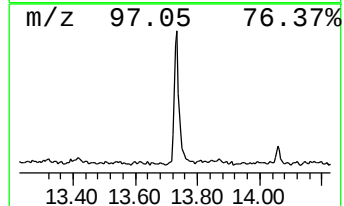
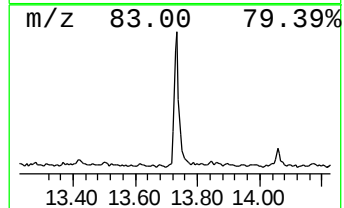
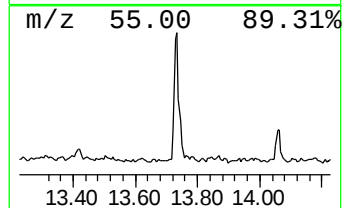
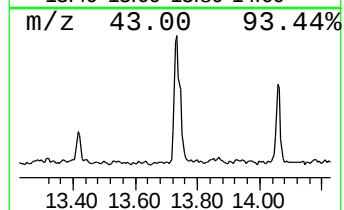
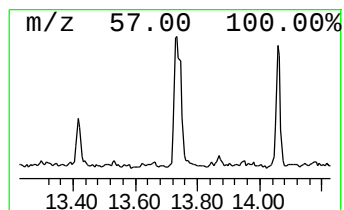
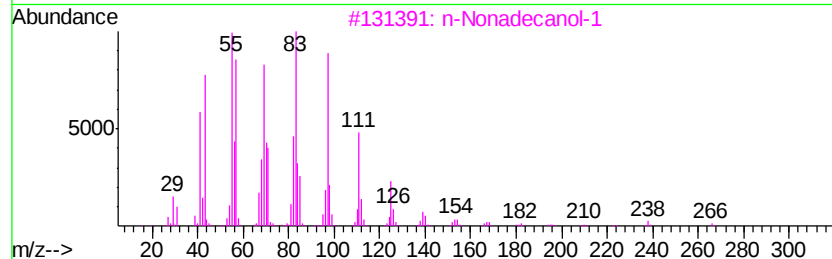
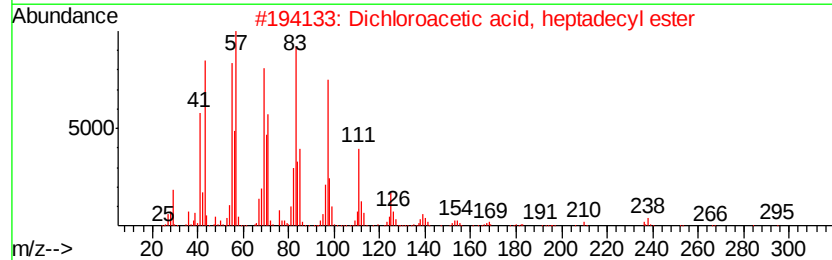
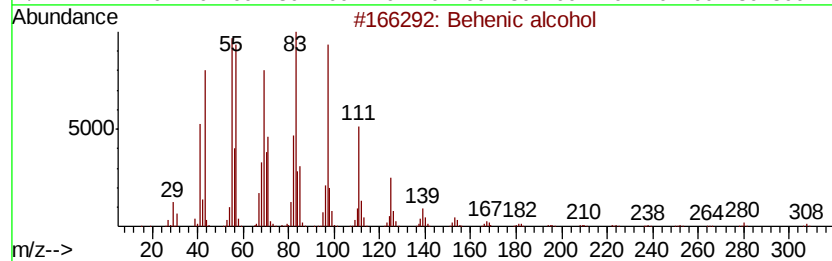
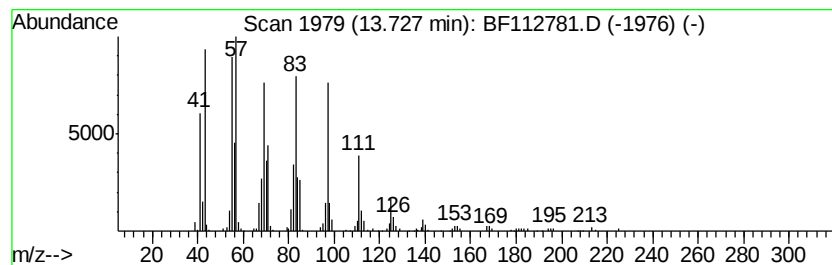
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Behenic alcohol Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD		R.T.	
13.73	9.37 ng	366958	Chrysene-d12		13.87	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Behenic alcohol	326	C22H46O	000661-19-8	95
2		Dichloroacetic acid, heptadecyl ...	366	C19H36Cl2O2	1000282-98-2	95
3		n-Nonadecanol-1	284	C19H40O	001454-84-8	95
4		Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	94
5		Bromoacetic acid, hexadecyl ester	362	C18H35BrO2	005454-48-8	94



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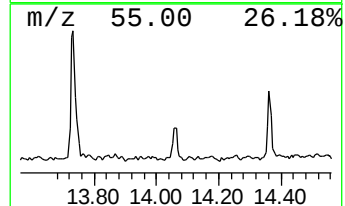
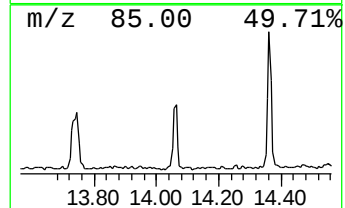
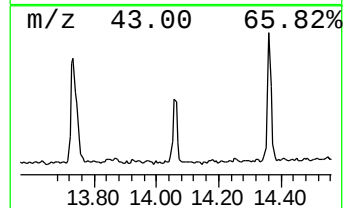
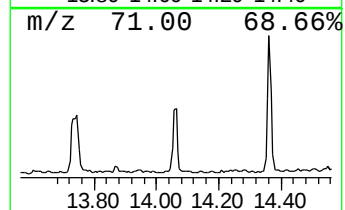
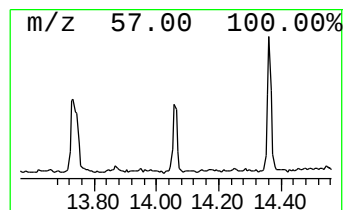
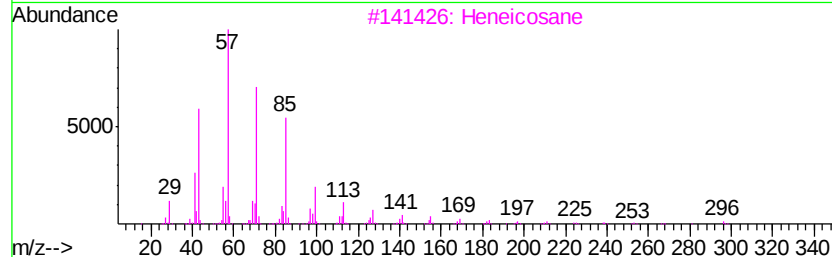
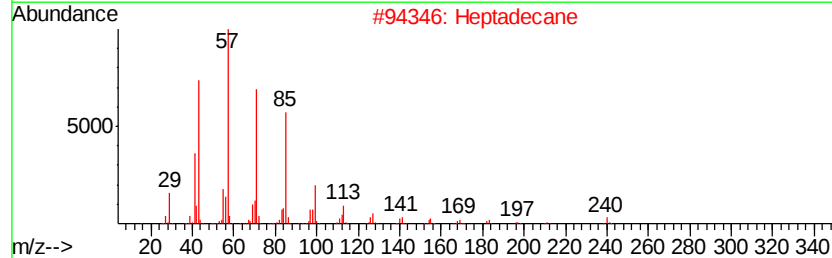
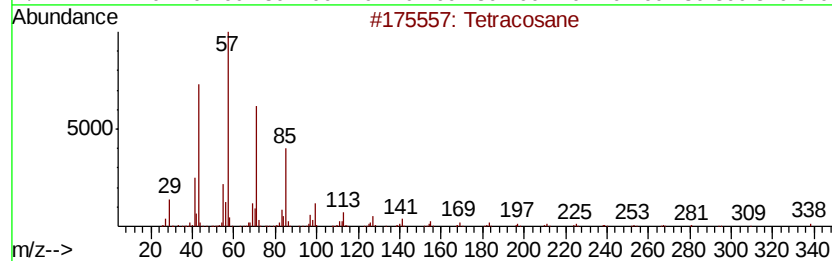
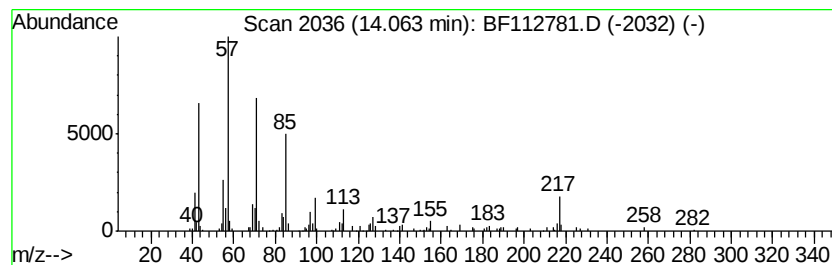
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Tetracosane Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.06	3.28 ng	128623	Chrysene-d12	13.87

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetracosane	338	C24H50	000646-31-1	76
2		Heptadecane	240	C17H36	000629-78-7	74
3		Heneicosane	296	C21H44	000629-94-7	74
4		Hentriacontane	437	C31H64	000630-04-6	74
5		Hexacosane	366	C26H54	000630-01-3	74



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF022819\
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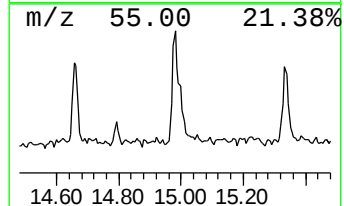
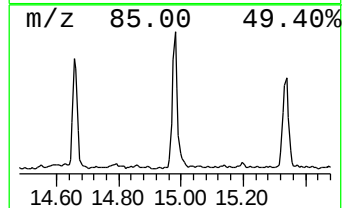
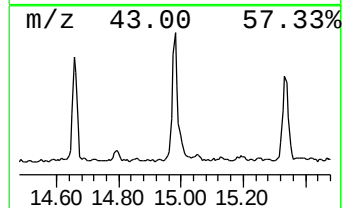
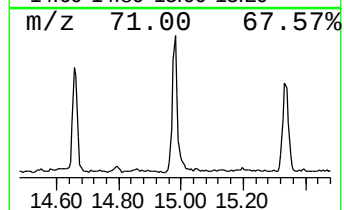
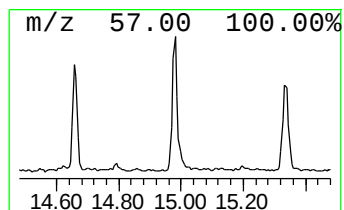
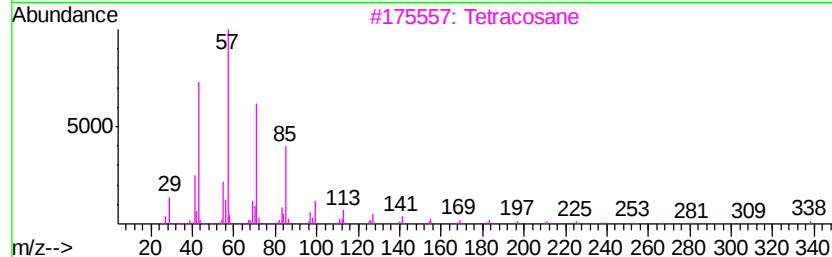
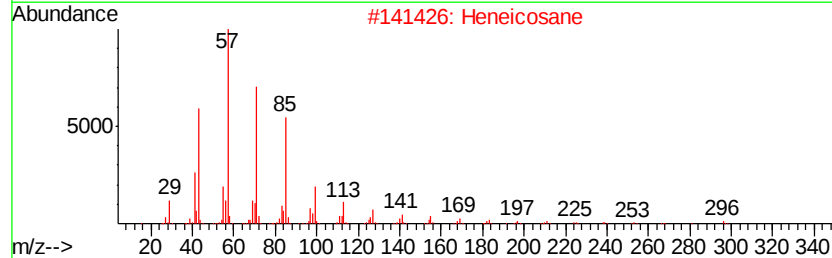
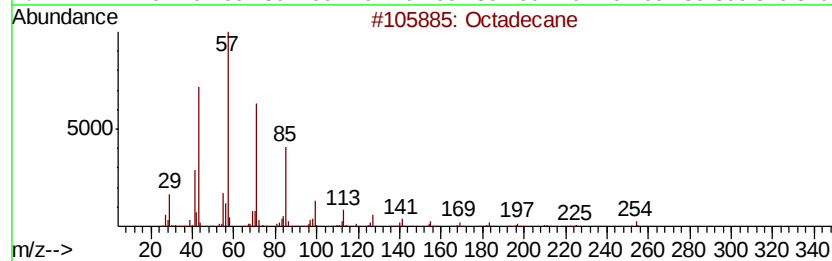
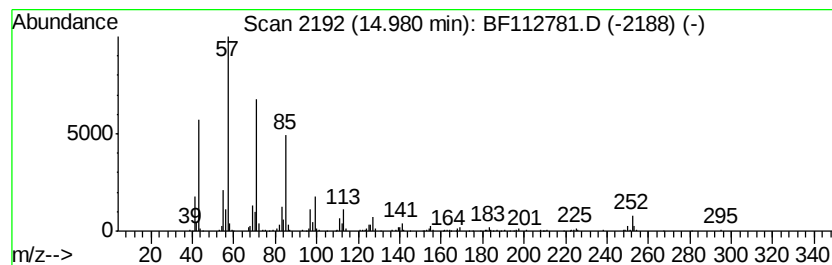
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Octadecane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.98	13.24 ng	423139	Perylene-d12	15.28

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecane	254	C18H38	000593-45-3	96
2		Heneicosane	296	C21H44	000629-94-7	91
3		Tetracosane	338	C24H50	000646-31-1	91
4		Eicosane, 7-hexyl-	366	C26H54	055333-99-8	90
5		Octadecane, 1-iodo-	380	C18H37I	000629-93-6	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF022819\
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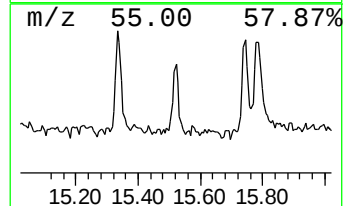
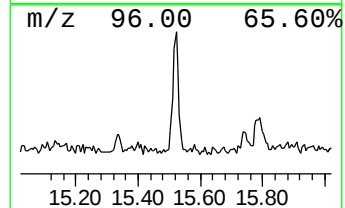
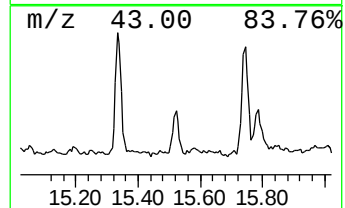
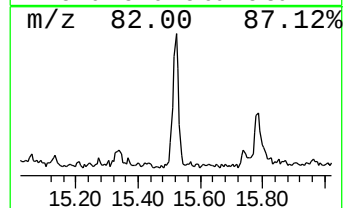
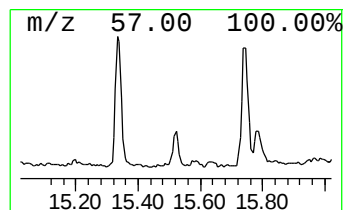
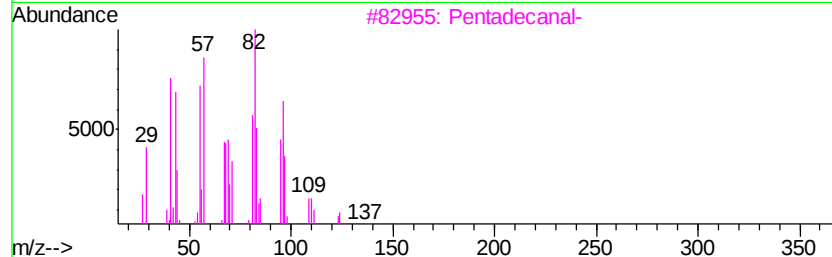
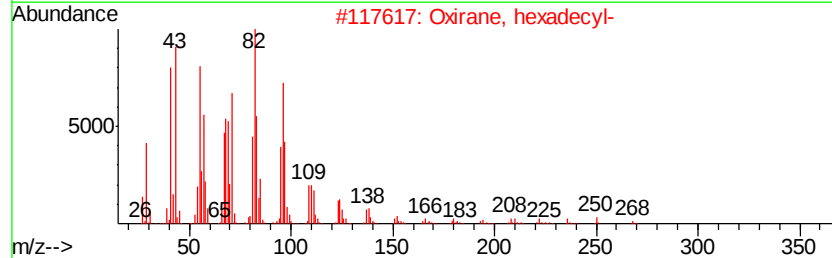
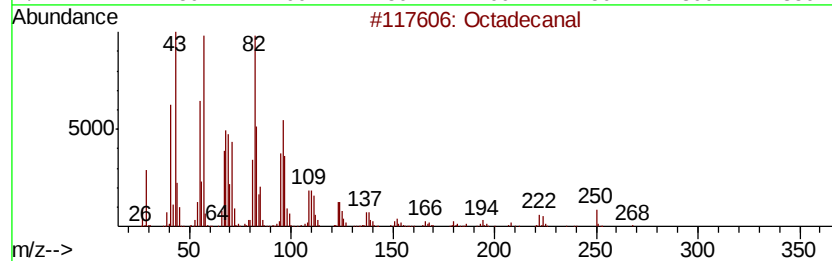
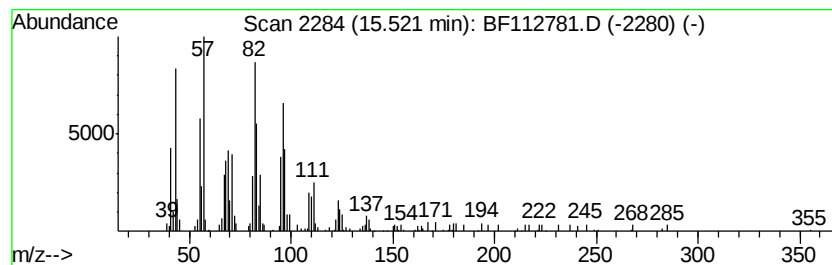
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 Octadecanal Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.52	4.97 ng	158742	Perylene-d12	15.28

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecanal	268	C18H36O	000638-66-4	90
2		Oxirane, hexadecyl-	268	C18H36O	007390-81-0	86
3		Pentadecanal-	226	C15H30O	002765-11-9	86
4		Oxirane, heptadecyl-	282	C19H38O	067860-04-2	86
5		Heptadecanal	254	C17H34O	1000376-70-0	86



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF022819\
Data File : BF112781.D
Acq On : 1 Mar 2019 00:27
Operator : JU/SJ
Sample : K1651-04
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
CON-327

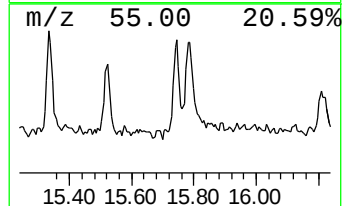
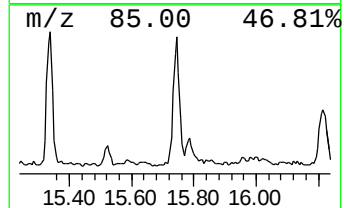
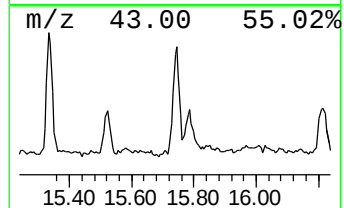
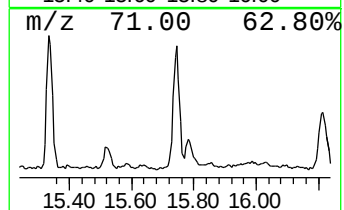
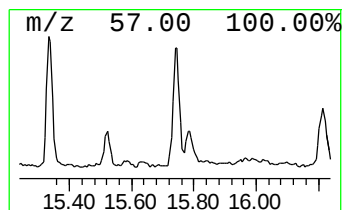
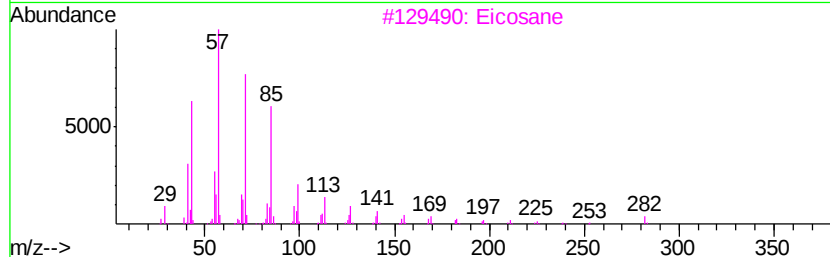
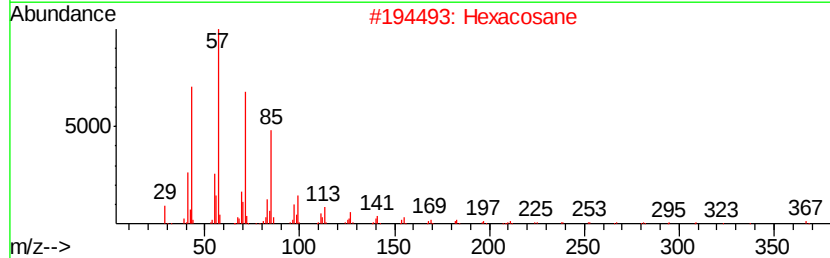
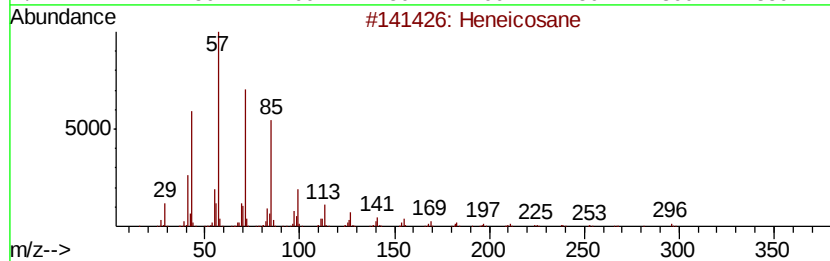
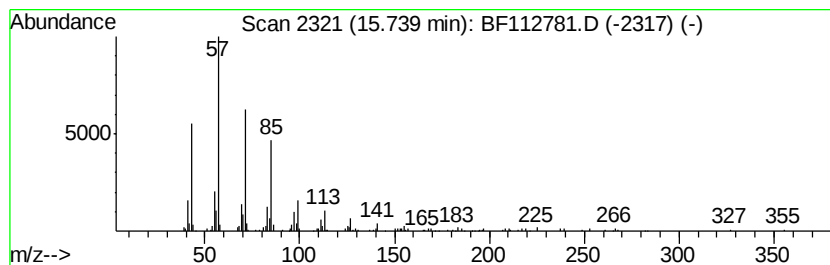
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 Heneicosane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.74	8.84 ng	282392	Perylene-d12	15.28

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosane	296	C21H44	000629-94-7	91
2		Hexacosane	366	C26H54	000630-01-3	90
3		Eicosane	282	C20H42	000112-95-8	90
4		Docosane	310	C22H46	000629-97-0	87
5		Decane, 3,8-dimethyl-	170	C12H26	017312-55-9	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF022819\
Data File : BF112781.D
Acq On : 1 Mar 2019 00:27
Operator : JU/SJ
Sample : K1651-04
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
CON-327

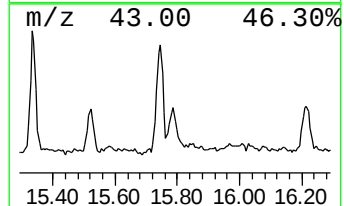
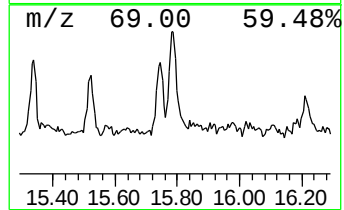
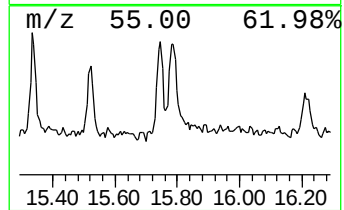
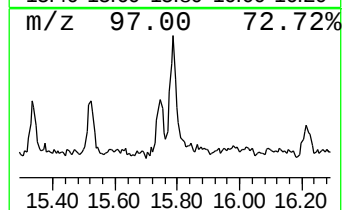
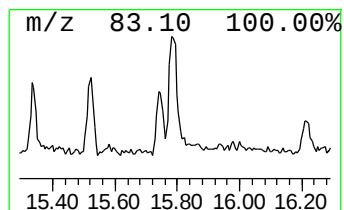
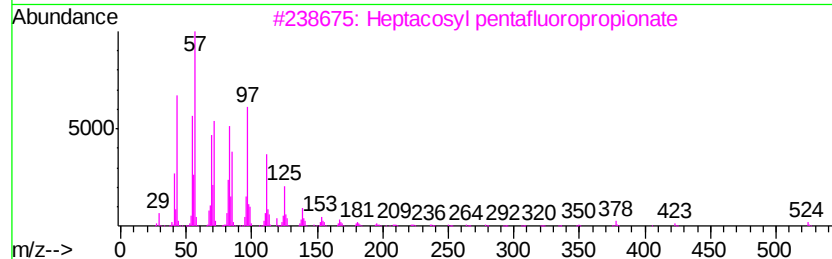
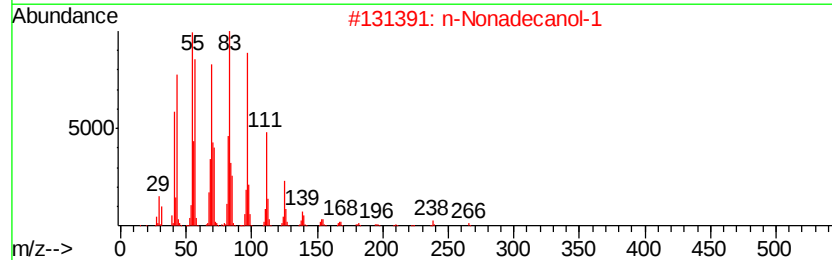
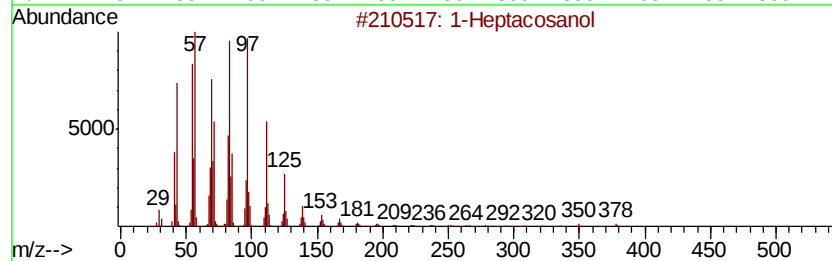
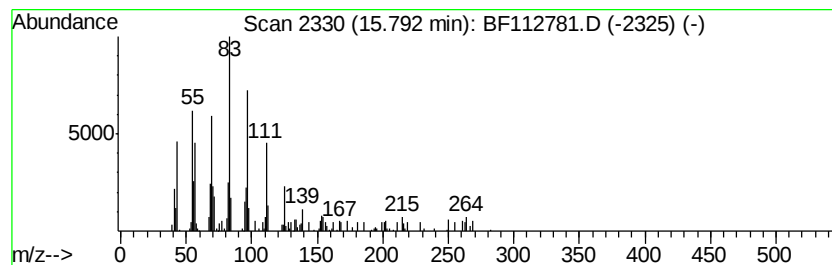
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 18 1-Heptacosanol Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.79	6.41 ng	204920	Perylene-d12	15.28

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Heptacosanol	396	C27H56O	002004-39-9	64
2		n-Nonadecanol-1	284	C19H40O	001454-84-8	64
3		Heptacosyl pentafluoropropionate	542	C30H55F5O2	1000351-81-6	53
4		Acetic acid n-octadecyl ester	312	C20H40O2	000822-23-1	53
5		1-Heneicosanol	312	C21H44O	015594-90-8	50



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF022819\
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Operator : JU/SJ
Sample : K1651-04
Misc :
ALS Vial : 17 Sample Multiplier: 1

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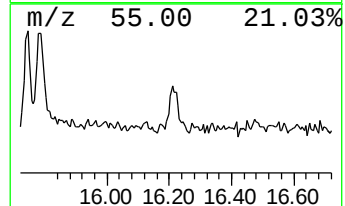
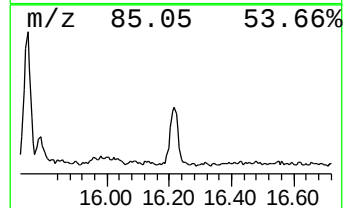
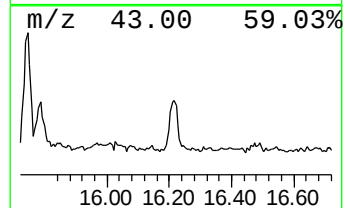
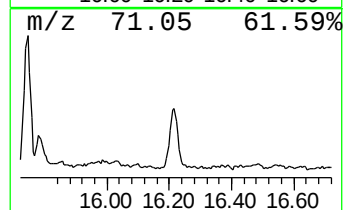
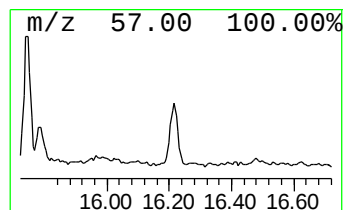
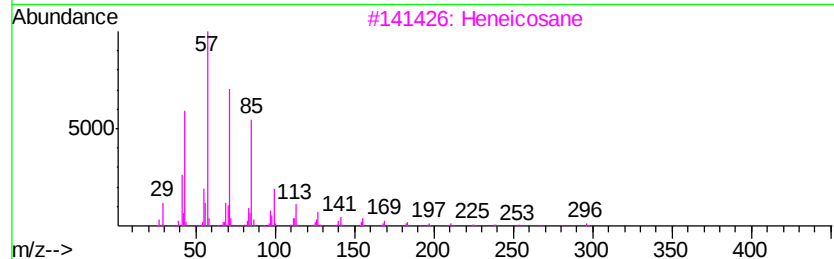
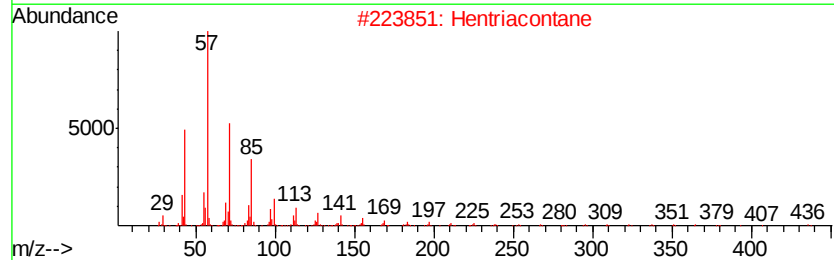
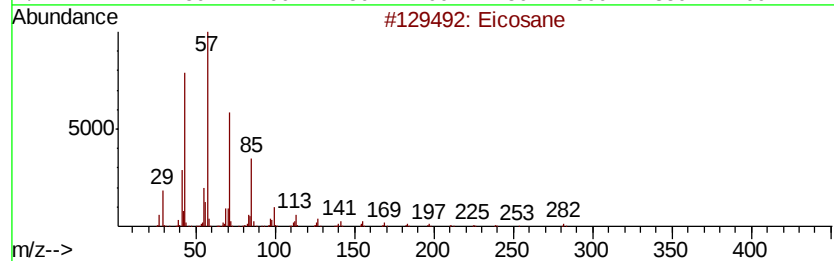
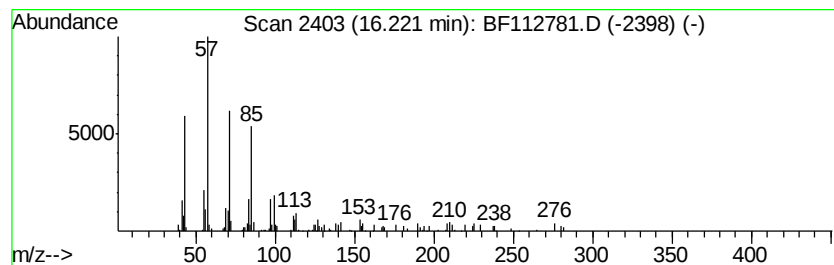
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 19 Eicosane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.22	4.98 ng	159130	Perylene-d12	15.28

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane	282	C20H42	000112-95-8	93
2		Hentriacontane	437	C31H64	000630-04-6	90
3		Heneicosane	296	C21H44	000629-94-7	87
4		Hexacosane	366	C26H54	000630-01-3	87
5		Pentacosane	352	C25H52	000629-99-2	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF022819\
 Data File : BF112781.D
 Acq On : 1 Mar 2019 00:27
 Operator : JU/SJ
 Sample : K1651-04
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CON-327

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF022719.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
3-Penten-2-one, (E)-	2.99	2.3	ng	63753	1	6.74	544920	20.0
4-Penten-2-one, 4...	3.27	5.0	ng	135260	1	6.74	544920	20.0
3-Penten-2-one, 4...	4.28	206.4	ng	5624530	1	6.74	544920	20.0
2-Pentanone, 4-hy...	4.94	211.8	ng	5771600	1	6.74	544920	20.0
unknown6.50	6.50	4.2	ng	115506	1	6.74	544920	20.0
2-Cyclohexen-1-on...	7.06	3.0	ng	83154	1	6.74	544920	20.0
2,4,7,9-Tetrameth...	9.22	4.2	ng	167418	3	9.77	796213	20.0
n-Nonadecanol-1	12.30	6.0	ng	240648	4	11.25	801358	20.0
Behenic alcohol	13.73	9.4	ng	366958	5	13.87	783493	20.0
Tetracosane	14.06	3.3	ng	128623	5	13.87	783493	20.0
Octadecane	14.98	13.2	ng	423139	6	15.28	639040	20.0
Octadecanal	15.52	5.0	ng	158742	6	15.28	639040	20.0
Heneicosane	15.74	8.8	ng	282392	6	15.28	639040	20.0
1-Heptacosanol	15.79	6.4	ng	204920	6	15.28	639040	20.0
Eicosane	16.22	5.0	ng	159130	6	15.28	639040	20.0