

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF110218\
 Data File : BF110437.D
 Acq On : 3 Nov 2018 2:54
 Operator : JU/SJ
 Sample : J5734-05
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 B-22(0-10)

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	2.293	30	35	49	rVB	347468	483229	21.90%	2.596%
2	4.098	337	342	351	rVB	65810	81286	3.68%	0.437%
3	5.204	526	530	539	rBV	61359	86133	3.90%	0.463%
4	5.586	590	595	598	rBV	2020238	1977459	89.60%	10.625%
5	6.586	759	765	779	rBV	1912617	2206864	100.00%	11.858%
6	6.728	784	789	792	rBV	2321761	2104651	95.37%	11.309%
7	6.957	824	828	833	rBV	615407	541785	24.55%	2.911%
8	7.110	850	854	858	rBV	1757786	1616476	73.25%	8.686%
9	7.522	919	924	934	rBV	1189134	1240372	56.21%	6.665%
10	8.239	1042	1046	1064	rBV	661834	655024	29.68%	3.520%
11	9.310	1224	1228	1237	rBV	2283577	2068732	93.74%	11.116%
12	9.704	1291	1295	1302	rBV	231281	219253	9.94%	1.178%
13	9.998	1341	1345	1359	rVB	685016	613357	27.79%	3.296%
14	10.786	1475	1479	1490	rBV	966046	967805	43.85%	5.200%
15	11.492	1595	1599	1607	rBV	518646	547505	24.81%	2.942%
16	11.992	1681	1684	1688	rBV2	36737	41037	1.86%	0.220%
17	12.551	1776	1779	1785	rVB7	14814	22442	1.02%	0.121%
18	12.704	1802	1805	1808	rBV	61940	59102	2.68%	0.318%
19	12.939	1839	1845	1850	rVB2	52318	76835	3.48%	0.413%
20	13.068	1863	1867	1871	rBV	1626499	1475720	66.87%	7.929%
21	13.274	1894	1902	1905	rBV2	22471	39055	1.77%	0.210%
22	13.539	1942	1947	1950	rBV	123600	132869	6.02%	0.714%
23	13.586	1952	1955	1958	rVV	75394	64479	2.92%	0.346%
24	13.621	1958	1961	1963	rVB2	23277	23325	1.06%	0.125%
25	13.951	2013	2017	2022	rBV	87392	102831	4.66%	0.553%
26	14.133	2044	2048	2052	rBV	409715	441045	19.99%	2.370%
27	14.262	2068	2070	2073	rVB	32216	31132	1.41%	0.167%
28	14.568	2119	2122	2125	rBV	47799	51780	2.35%	0.278%
29	14.739	2148	2151	2155	rVB4	29317	31430	1.42%	0.169%
30	14.886	2173	2176	2179	rVB	64439	59306	2.69%	0.319%
31	15.239	2233	2236	2240	rVB	73208	82906	3.76%	0.445%
32	15.662	2300	2308	2315	rVB2	212245	353559	16.02%	1.900%
33	16.092	2378	2381	2388	rVB2	51780	71742	3.25%	0.385%
34	16.621	2468	2471	2476	rVB3	27626	40660	1.84%	0.218%

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Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
B-22(0-10)

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

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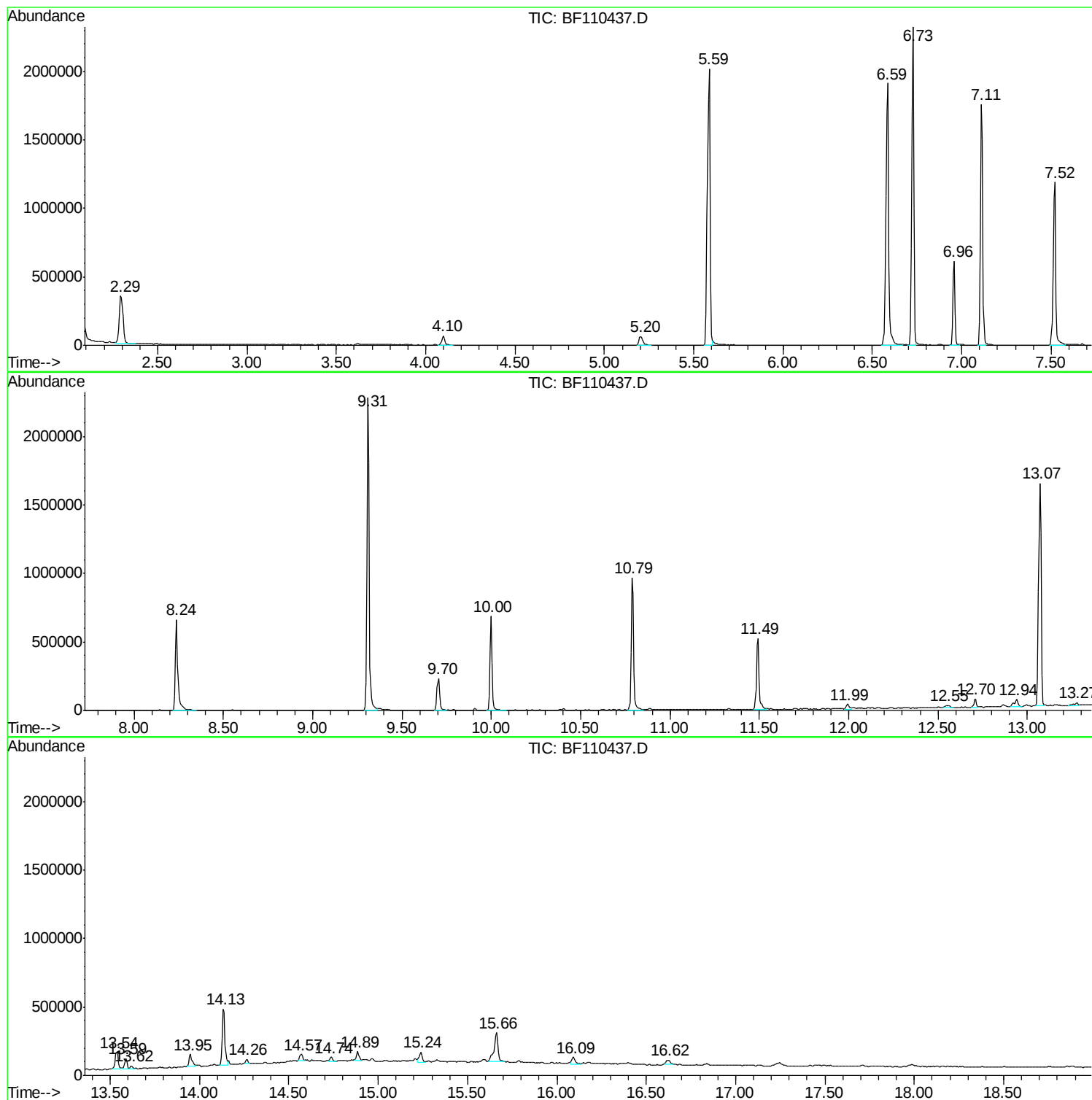
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ClientSampled :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF102318.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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ClientSampleId :
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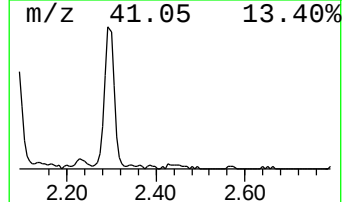
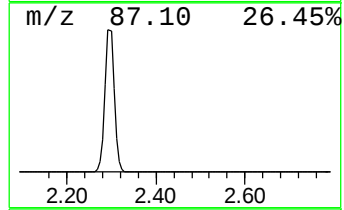
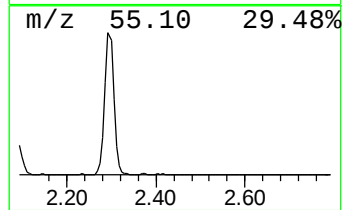
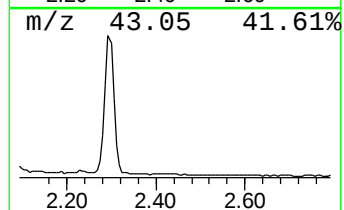
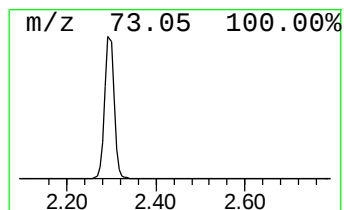
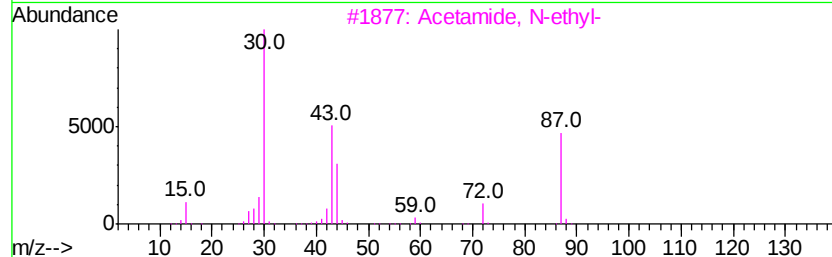
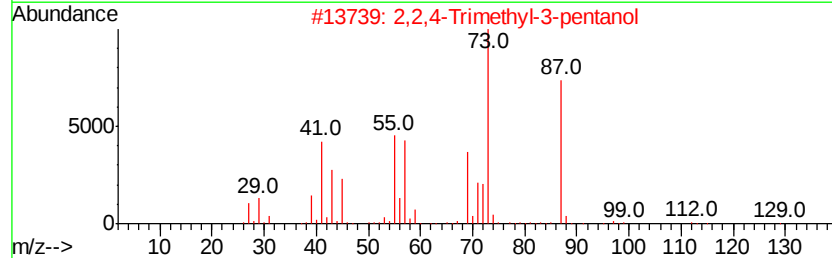
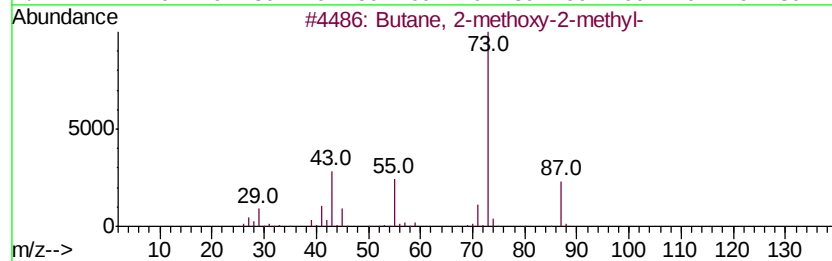
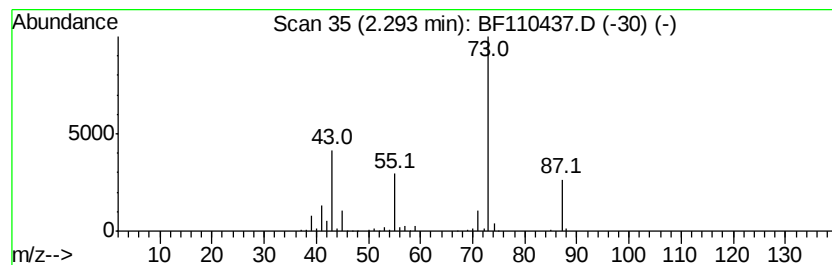
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF102318.M
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.29	17.84 ng	483229	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		Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	27
4		Silane, tetramethyl-	88	C4H12Si	000075-76-3	12
5		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	12



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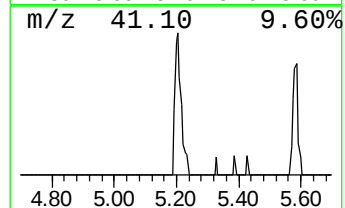
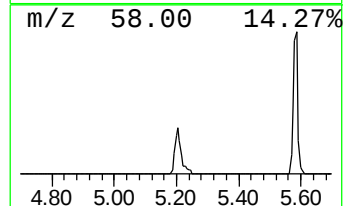
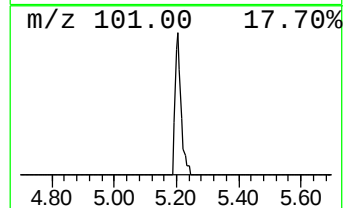
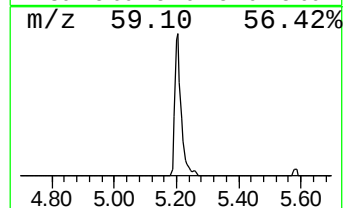
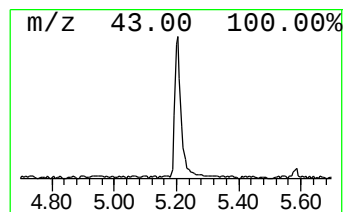
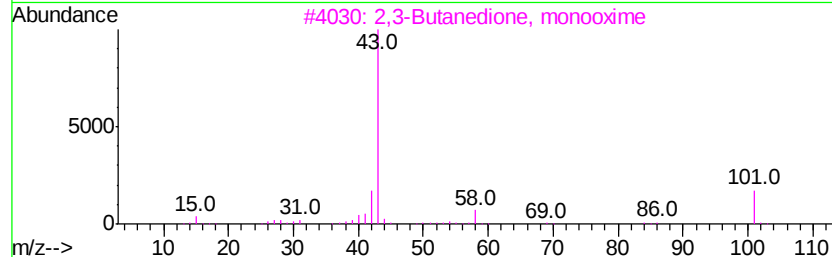
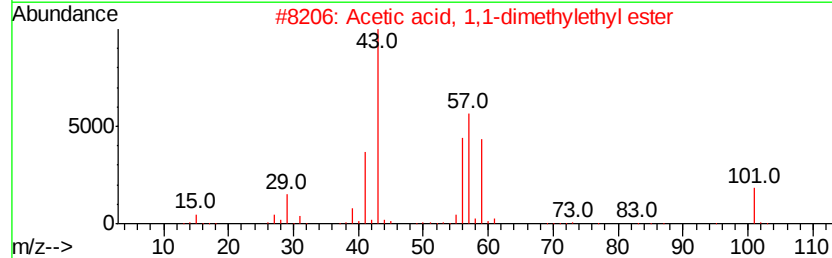
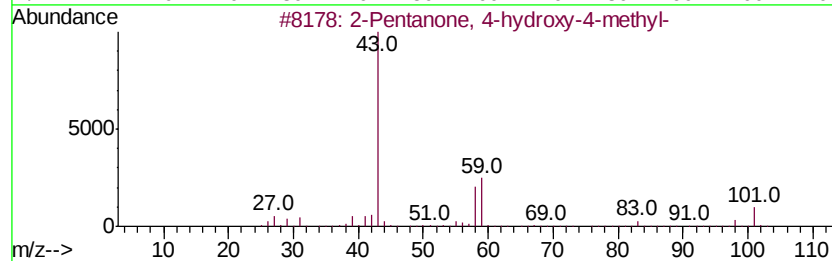
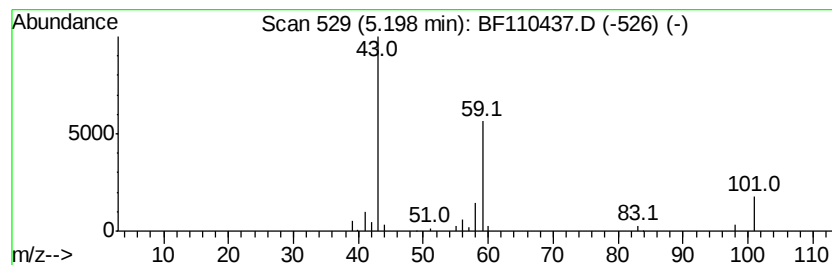
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.20	3.18 ng	86133	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	59
2		Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	38
3		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
4		Ethanone, 1-(2-methylcyclopropyl)-	98	C6H10O	000930-56-3	9
5		3-Heptanol, 4-methyl-	130	C8H18O	014979-39-6	9



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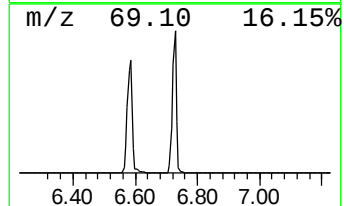
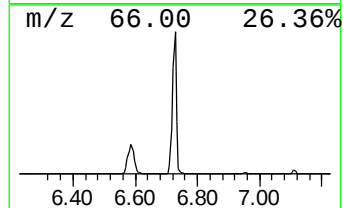
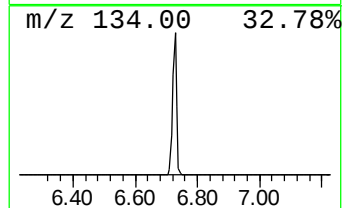
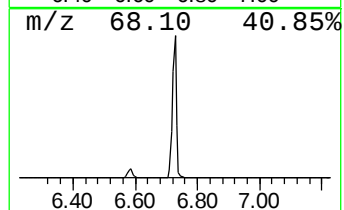
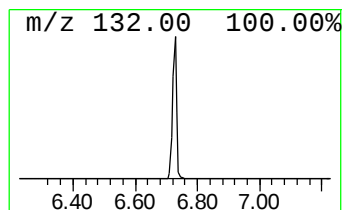
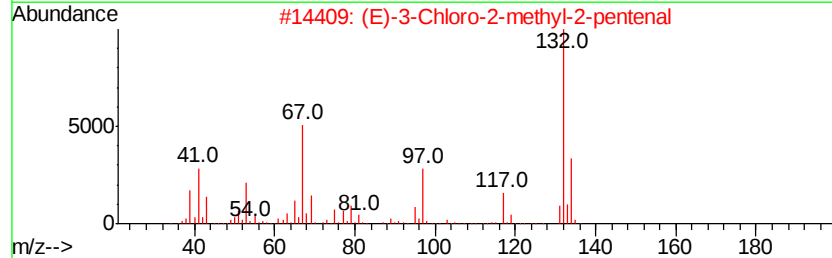
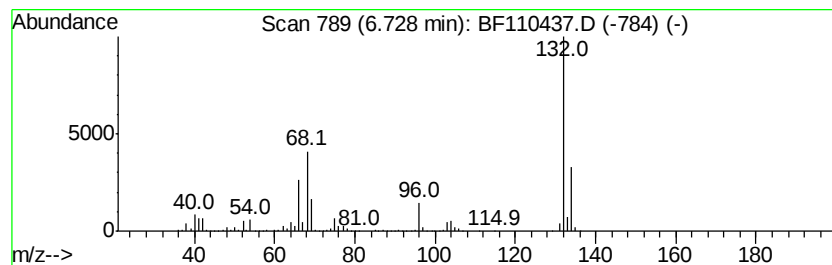
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 unknown6.73 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.73	77.69 ng	2104650	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		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	12
3		Tranvlcvpromine-propionyl	189	C12H15NO	1000123-86-3	12
4		Xenon	132	Xe	007440-63-3	9
5		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9



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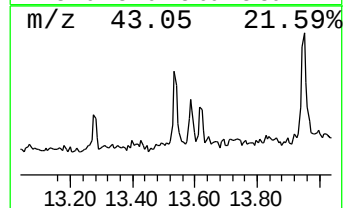
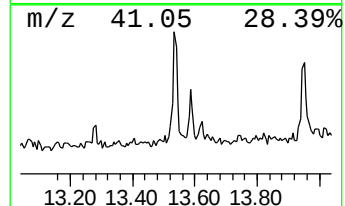
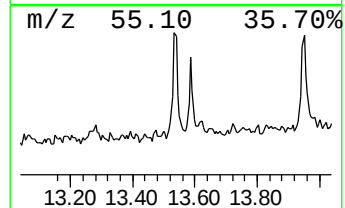
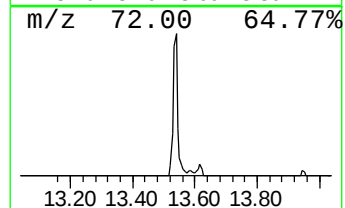
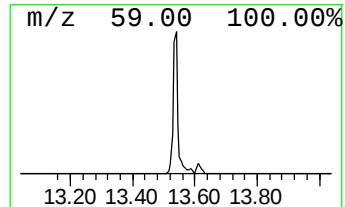
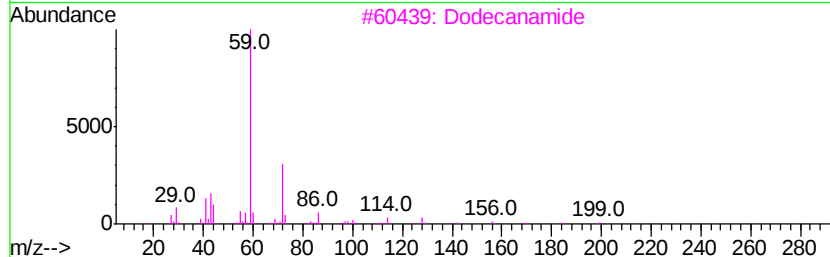
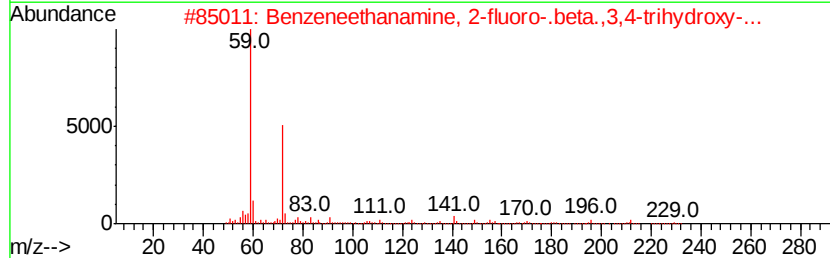
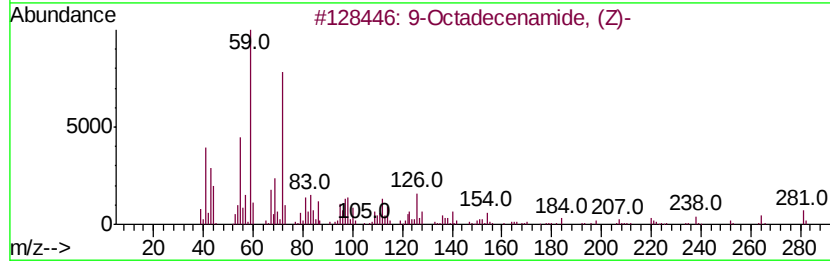
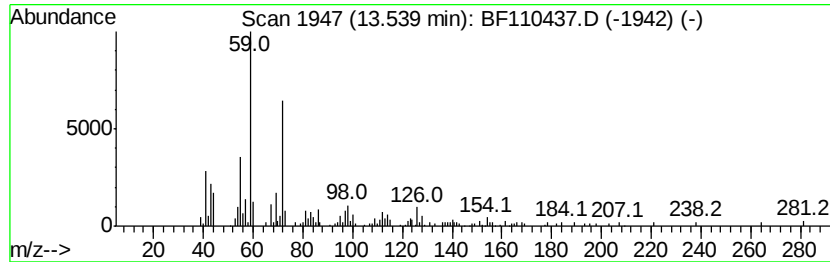
Instrument :
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 9-Octadecenamide, (Z)- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD		R.T.
13.54	6.03 ng	132869	Chrysene-d12		14.13
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	95
2	Benzeneethanamine, 2-fluoro-.bet...	229	C11H16FN03	061338-98-5	59
3	Dodecanamide	199	C12H25NO	001120-16-7	53
4	Octanamide	143	C8H17NO	000629-01-6	53
5	Acetic acid, cyanohydroxyimino-,...	128	C4H4N2O3	061295-92-9	43



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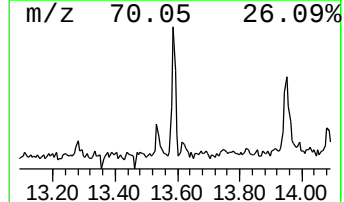
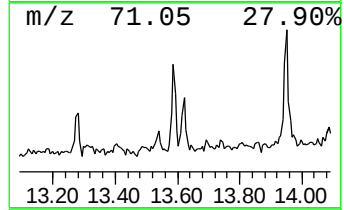
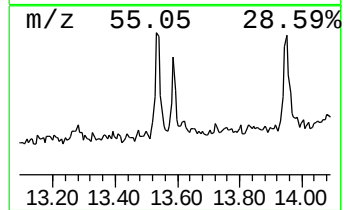
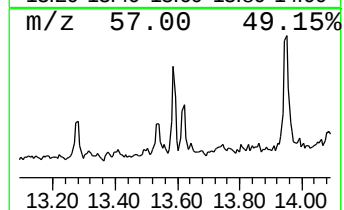
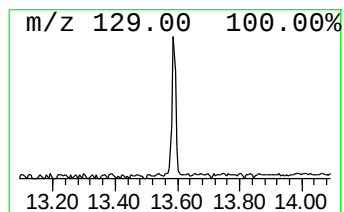
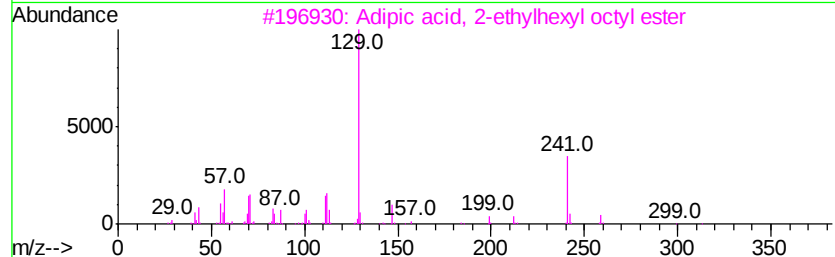
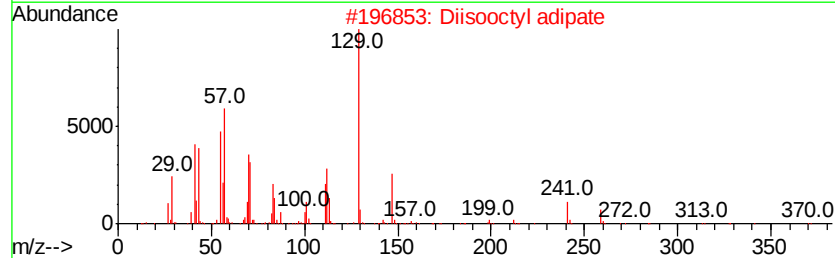
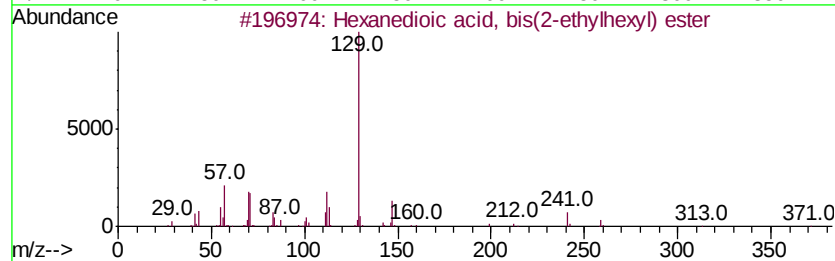
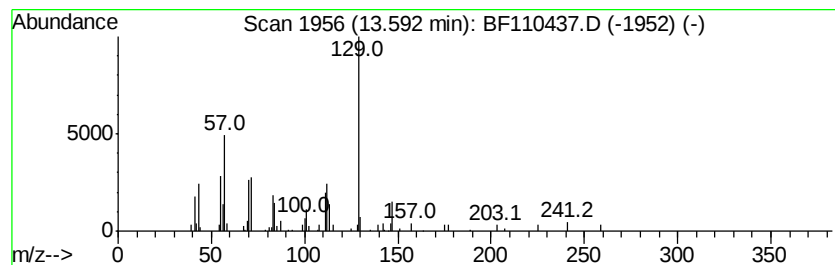
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 Hexanedioic acid, bis(2-eth... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.59	2.92 ng	64479	Chrysene-d12	14.13

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexanedioic acid, bis(2-ethylhex...	370	C22H42O4	000103-23-1	91
2		Diisooctyl adipate	370	C22H42O4	001330-86-5	86
3		Adipic acid, 2-ethylhexyl octyl ...	370	C22H42O4	1000324-48-6	83
4		Adipic acid, 2-ethylhexyl isobut...	314	C18H34O4	1000324-48-0	72
5		Hexanedioic acid, mono(2-ethylhe...	258	C14H26O4	004337-65-9	58



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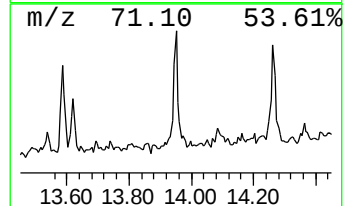
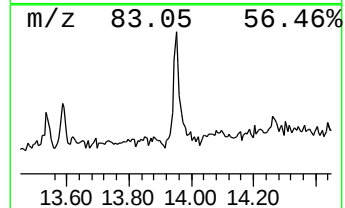
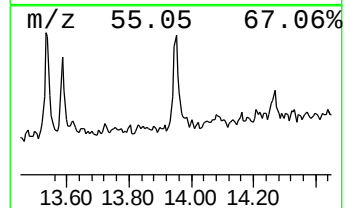
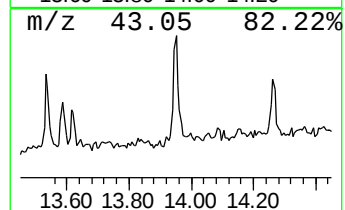
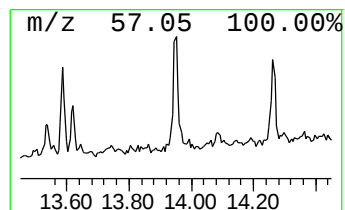
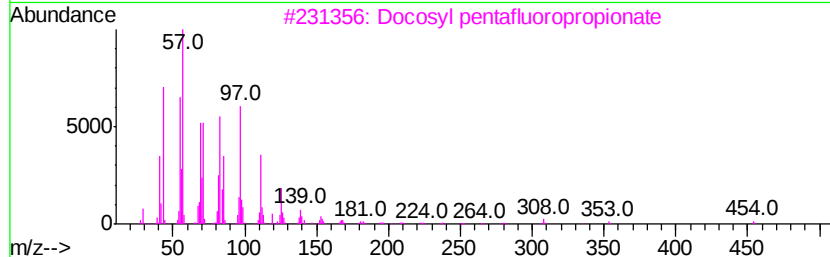
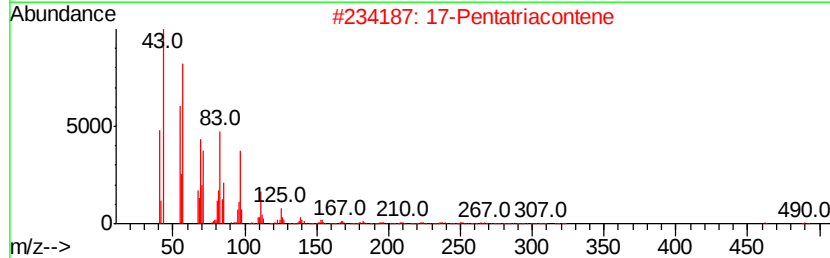
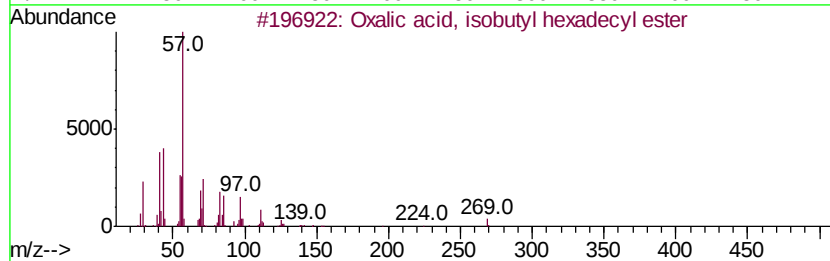
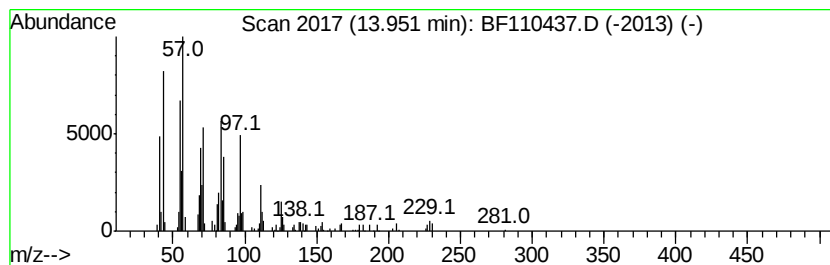
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Oxalic acid, isobutyl hexad... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.95	4.66 ng	102831	Chrysene-d12	14.13

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Oxalic acid, isobutyl hexadecyl ...	370	C22H42O4	1000309-38-1	91
2			17-Pentatriacontene	491	C35H70	006971-40-0	90
3			Docosyl pentafluoropropionate	472	C25H45F5O2	1000351-80-9	87
4			Dotriacontyl heptafluorobutyrate	662	C36H65F7O2	1000351-84-2	87
5			Ethanol, 2-(tetradecyloxy)-	258	C16H34O2	002136-70-1	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF110218\
Data File : BF110437.D
Acq On : 3 Nov 2018 2:54
Operator : JU/SJ
Sample : J5734-05
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
B-22(0-10)

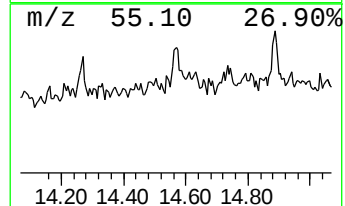
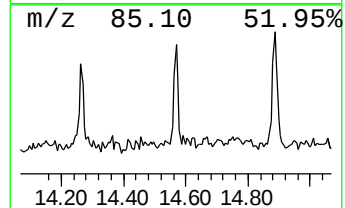
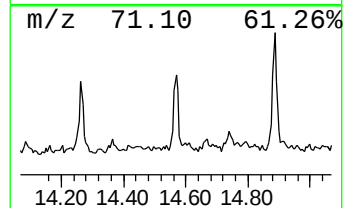
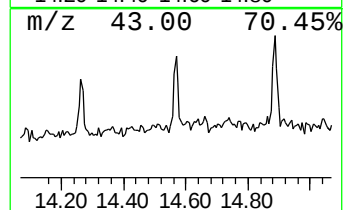
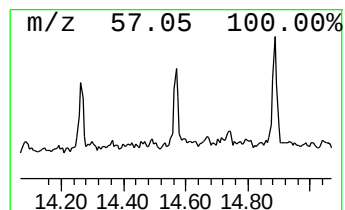
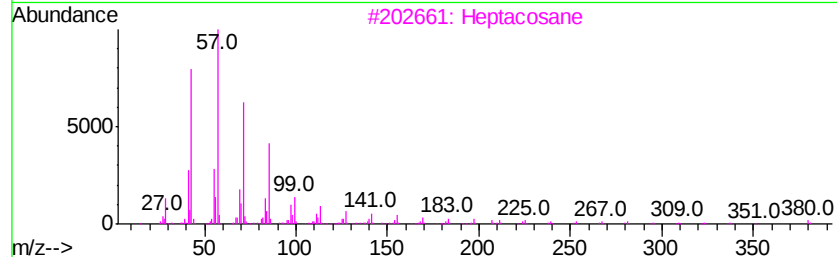
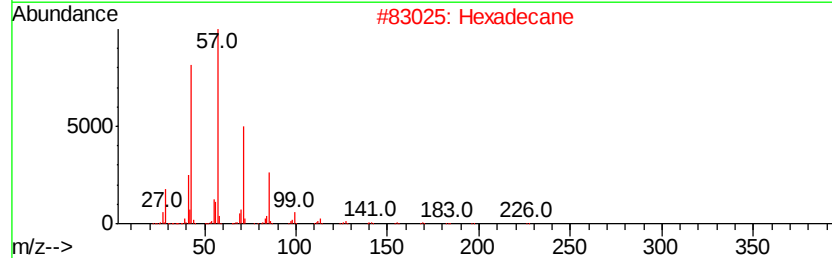
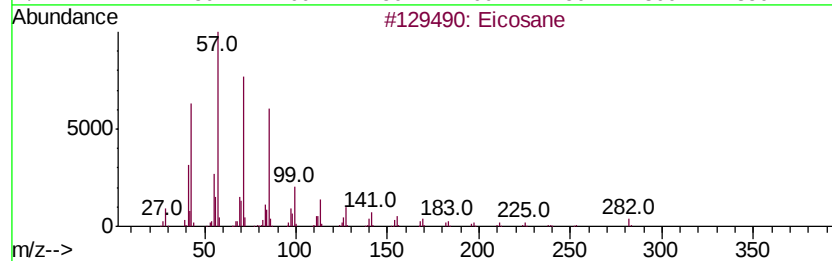
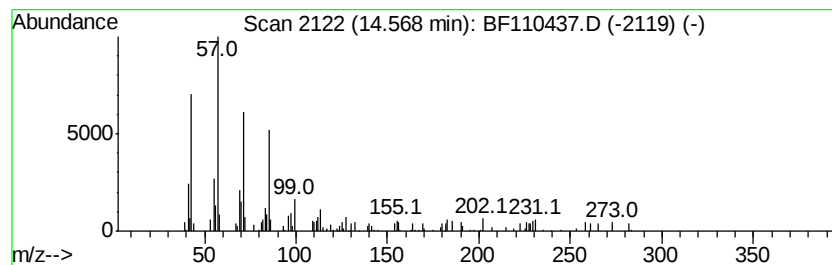
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Eicosane Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.57	2.35 ng	51780	Chrysene-d12	14.13

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane	282	C20H42	000112-95-8	93
2		Hexadecane	226	C16H34	000544-76-3	86
3		Heptacosane	380	C27H56	000593-49-7	62
4		Tridecanol, 2-ethyl-2-methyl-	242	C16H34O	1000115-66-1	62
5		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	62



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF110218\
Data File : BF110437.D
Acq On : 3 Nov 2018 2:54
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Sample : J5734-05
Misc :
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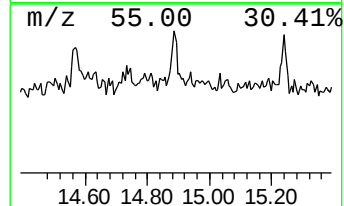
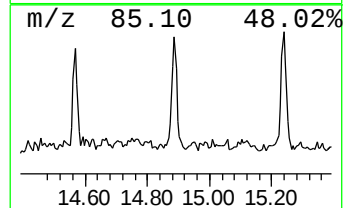
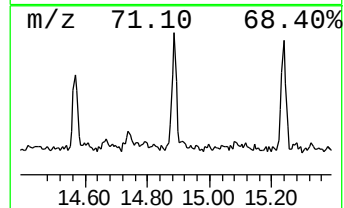
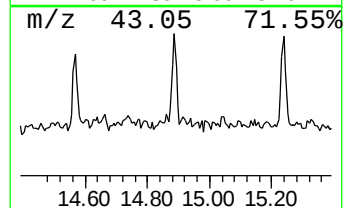
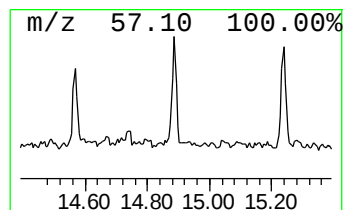
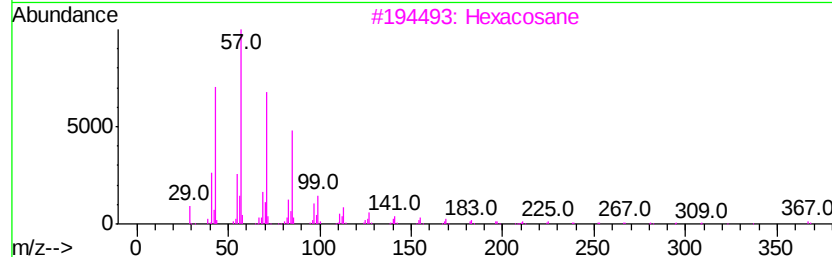
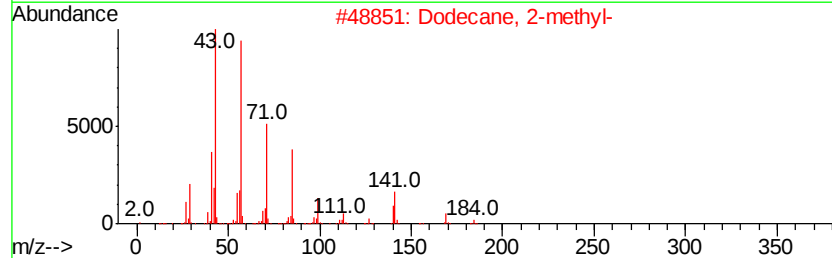
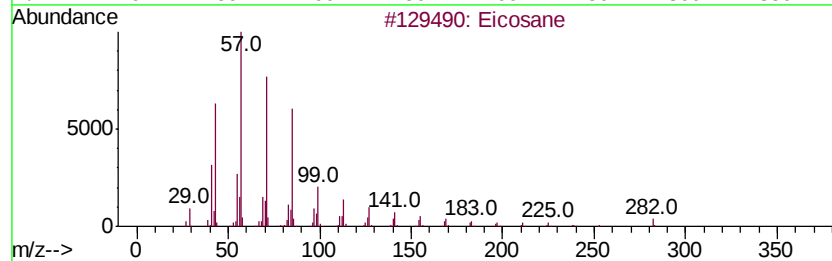
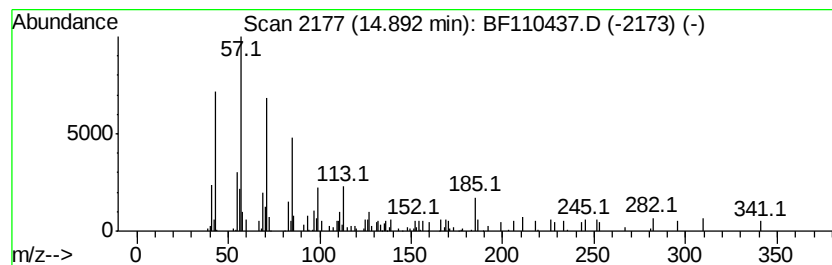
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Dodecane, 2-methyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.89	2.69 ng	59306	Chrysene-d12	14.13
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Eicosane	282 C20H42	000112-95-8	64
2	Dodecane, 2-methyl-	184 C13H28	001560-97-0	58
3	Hexacosane	366 C26H54	000630-01-3	58
4	Hexadecane	226 C16H34	000544-76-3	58
5	Heptadecane	240 C17H36	000629-78-7	58



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF110218\
Data File : BF110437.D
Acq On : 3 Nov 2018 2:54
Operator : JU/SJ
Sample : J5734-05
Misc :
ALS Vial : 26 Sample Multiplier: 1

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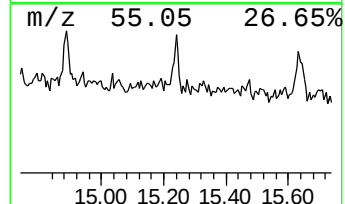
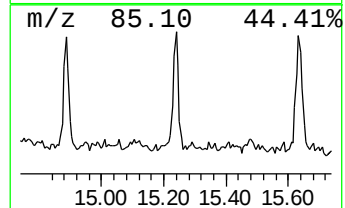
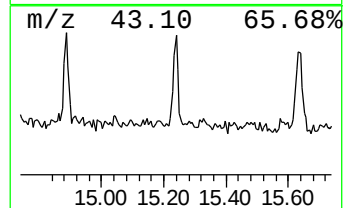
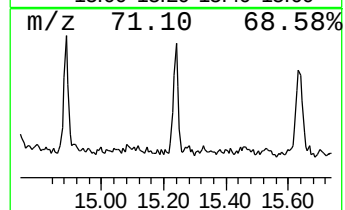
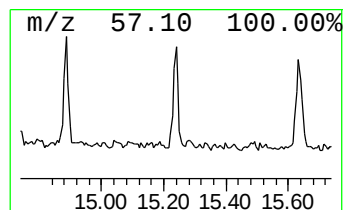
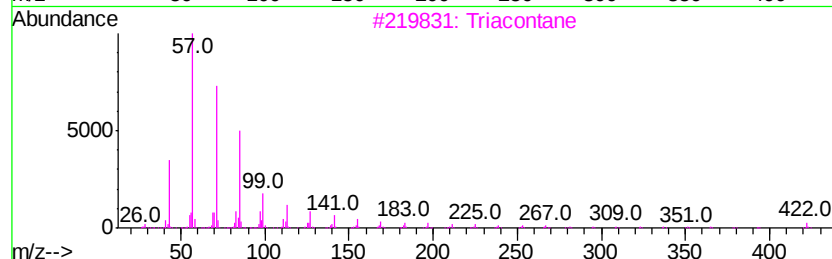
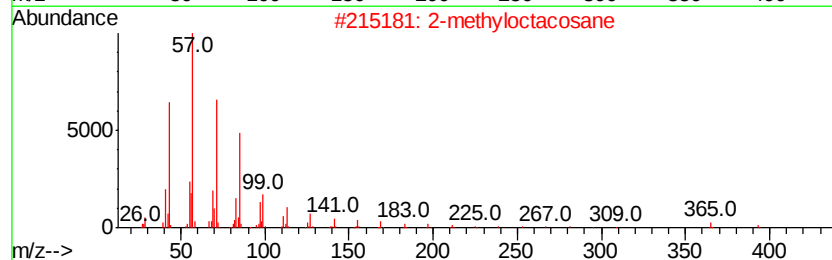
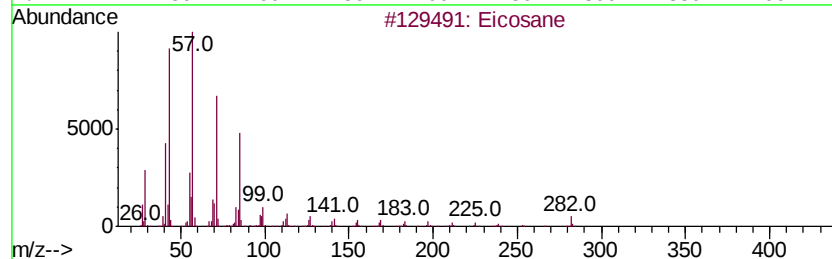
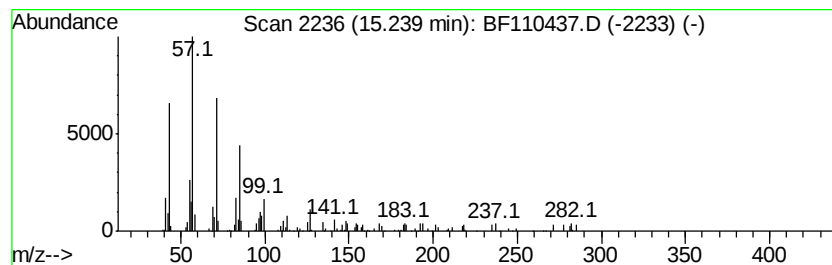
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 2-methyloctacosane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.24	4.69 ng	82906	Perylene-d12	15.66

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Eicosane	282	C20H42	000112-95-8	90
2			2-methyloctacosane	408	C29H60	1000376-72-8	83
3			Triacontane	422	C30H62	000638-68-6	80
4			Tridecane, 6-propyl-	226	C16H34	055045-10-8	80
5			Heneicosane, 11-(1-ethylpropyl)-	366	C26H54	055282-11-6	80



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF110218\
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Sample : J5734-05
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ALS Vial : 26 Sample Multiplier: 1

Instrument :
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ClientSampleId :
B-22(0-10)

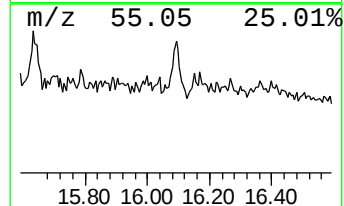
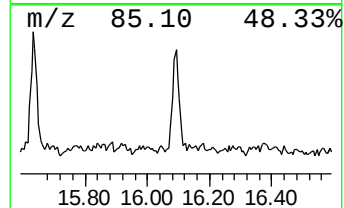
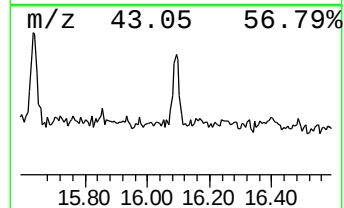
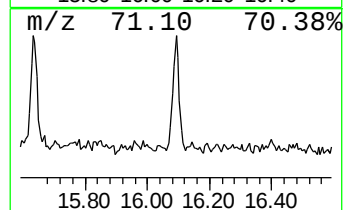
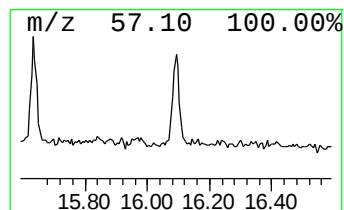
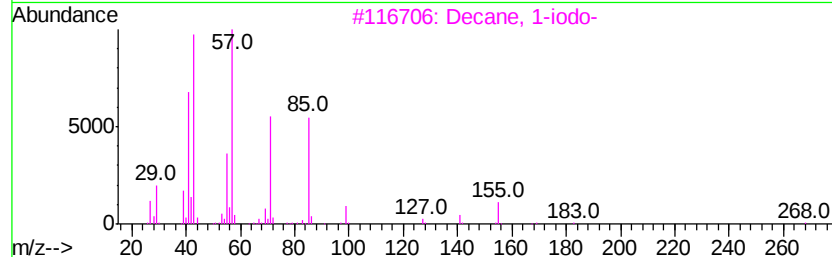
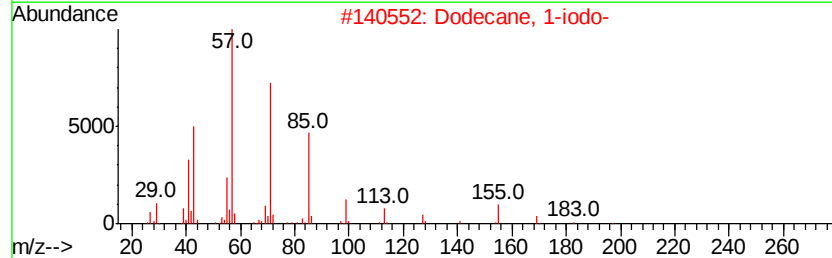
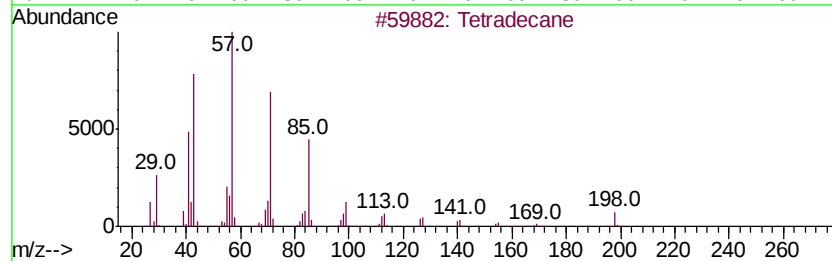
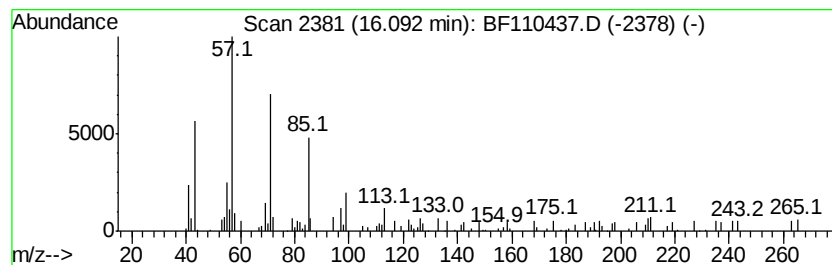
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Tetradecane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.09	4.06 ng	71742	Perylene-d12	15.66

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetradecane	198	C14H30	000629-59-4	64
2		Dodecane, 1-iodo-	296	C12H25I	004292-19-7	50
3		Decane, 1-iodo-	268	C10H21I	002050-77-3	47
4		2,5-Heptanedione, 3,3,6-trimethyl-	170	C10H18O2	051513-40-7	47
5		5-Tridecanone	198	C13H26O	030692-16-1	38



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF110218\
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ALS Vial : 26 Sample Multiplier: 1

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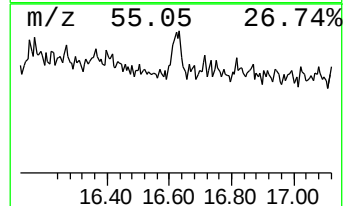
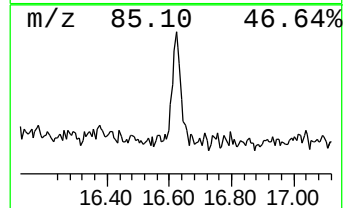
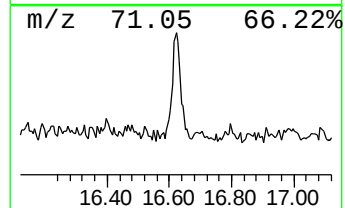
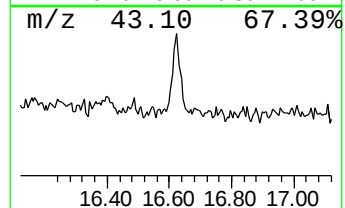
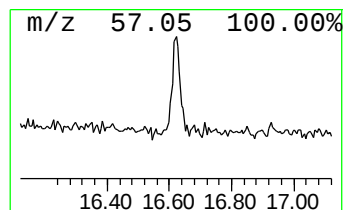
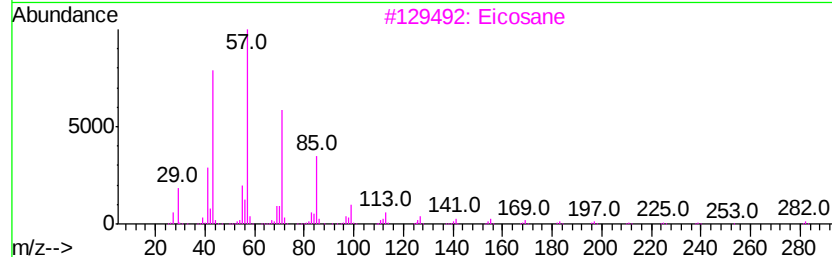
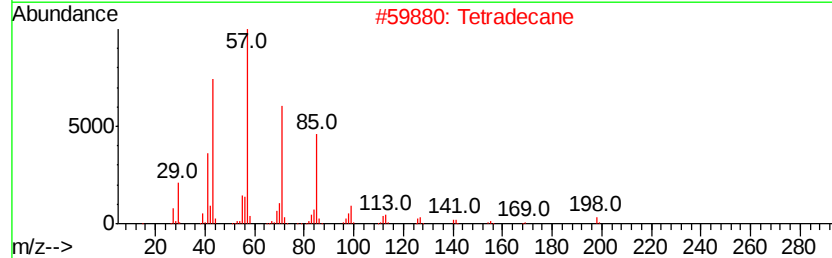
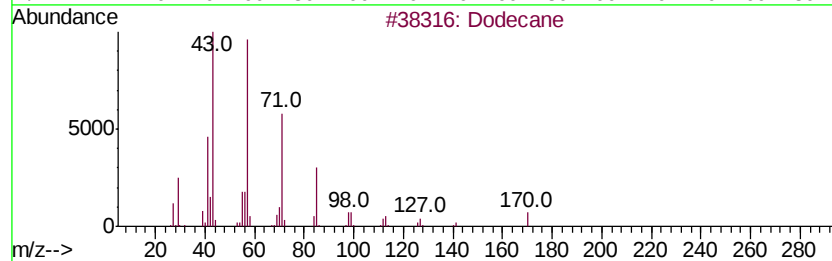
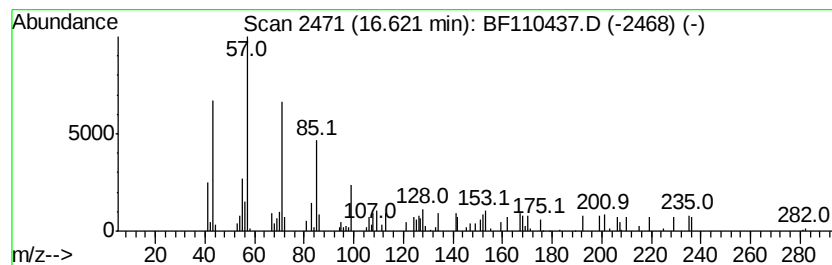
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Dodecane Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.62	2.30 ng	40660	Perylene-d12	15.66

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Dodecane	170	C12H26	000112-40-3	52
2		Tetradecane	198	C14H30	000629-59-4	50
3		Eicosane	282	C20H42	000112-95-8	50
4		2,5-Heptanedione, 3,3,6-trimethyl-	170	C10H18O2	051513-40-7	47
5		Dodecane, 5-methyl-	184	C13H28	017453-93-9	43



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF110218\
Data File : BF110437.D
Acq On : 3 Nov 2018 2:54
Operator : JU/SJ
Sample : J5734-05
Misc :
ALS Vial : 26 Sample Multiplier: 1

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

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-methoxy...	2.29	17.8	ng	483229	1	6.96	541785	20.0
2-Pentanone, 4-hy...	5.20	3.2	ng	86133	1	6.96	541785	20.0
unknown6.73	6.73	77.7	ng	2104650	1	6.96	541785	20.0
9-Octadecenamide,...	13.54	6.0	ng	132869	5	14.13	441045	20.0
Hexanedioic acid,...	13.59	2.9	ng	64479	5	14.13	441045	20.0
Oxalic acid, isob...	13.95	4.7	ng	102831	5	14.13	441045	20.0
Eicosane	14.57	2.4	ng	51780	5	14.13	441045	20.0
Dodecane, 2-methyl-	14.89	2.7	ng	59306	5	14.13	441045	20.0
2-methyloctacosane	15.24	4.7	ng	82906	6	15.66	353559	20.0
Tetradecane	16.09	4.1	ng	71742	6	15.66	353559	20.0
Dodecane	16.62	2.3	ng	40660	6	15.66	353559	20.0