

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF111919\
 Data File : BF117879.D
 Acq On : 20 Nov 2019 3:27
 Operator : JU
 Sample : K5908-02
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 B-8-1-(6-12)

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-BF111319.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	2.199	13	19	30	rVB	674287	854650	21.65%	2.378%
2	5.128	512	517	524	rVB	559632	635618	16.10%	1.768%
3	5.528	580	585	588	rBV	3413384	3158655	80.03%	8.788%
4	6.534	751	756	764	rBV	3200325	3491107	88.45%	9.713%
5	6.681	777	781	785	rBV	3597109	3434351	87.02%	9.555%
6	6.916	817	821	824	rBV	1356278	1049236	26.58%	2.919%
7	7.075	844	848	851	rBV	3008843	2711826	68.71%	7.545%
8	7.475	907	916	919	rBV	2307406	2066363	52.36%	5.749%
9	8.204	1033	1040	1043	rBV	1322571	1169643	29.64%	3.254%
10	9.016	1175	1178	1181	rBV	60991	49692	1.26%	0.138%
11	9.281	1218	1223	1226	rBV	4716100	3946797	100.00%	10.981%
12	9.392	1238	1242	1246	rVB2	134446	138225	3.50%	0.385%
13	9.622	1278	1281	1283	rBV	55300	50187	1.27%	0.140%
14	9.675	1287	1290	1294	rVB	714909	584998	14.82%	1.628%
15	9.963	1333	1339	1343	rBV	1479508	1320002	33.44%	3.672%
16	10.375	1405	1409	1411	rBV2	38420	52941	1.34%	0.147%
17	10.757	1469	1474	1477	rBV	2057187	1952824	49.48%	5.433%
18	10.880	1491	1495	1500	rVB	87222	109354	2.77%	0.304%
19	11.457	1589	1593	1596	rBV	1374105	1159844	29.39%	3.227%
20	11.480	1596	1597	1599	rVB	100567	50600	1.28%	0.141%
21	11.980	1676	1682	1686	rBV2	510419	552908	14.01%	1.538%
22	12.516	1769	1773	1775	rBV3	77494	116079	2.94%	0.323%
23	12.545	1776	1778	1781	rVB2	75401	71745	1.82%	0.200%
24	12.674	1797	1800	1803	rBV	115628	127489	3.23%	0.355%
25	12.769	1812	1816	1825	rBV	373764	461338	11.69%	1.284%
26	12.904	1834	1839	1840	rBV2	91576	93400	2.37%	0.260%
27	13.045	1859	1863	1867	rBV	3615729	3210792	81.35%	8.933%
28	13.274	1900	1902	1907	rBV3	70735	94599	2.40%	0.263%
29	13.604	1955	1958	1959	rBV2	89095	74843	1.90%	0.208%
30	13.739	1979	1981	1985	rBV	46817	49057	1.24%	0.136%
31	13.951	2014	2017	2018	rBV	93908	92893	2.35%	0.258%
32	13.968	2018	2020	2028	rVB2	128707	252542	6.40%	0.703%
33	14.104	2039	2043	2046	rVB	1138383	980623	24.85%	2.728%
34	14.263	2068	2070	2073	rVB	123742	121220	3.07%	0.337%

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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

Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF111319.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	14.568	2120	2122	2127	rVB	151765	148437	3.76%	0.413%
36	14.892	2174	2177	2181	rVB	126698	127342	3.23%	0.354%
37	15.245	2234	2237	2240	rVB	139370	142181	3.60%	0.396%
38	15.610	2295	2299	2303	rBV	631619	802915	20.34%	2.234%
39	16.104	2380	2383	2389	rVB	114287	190909	4.84%	0.531%
40	17.345	2589	2594	2601	rVB5	115124	245060	6.21%	0.682%

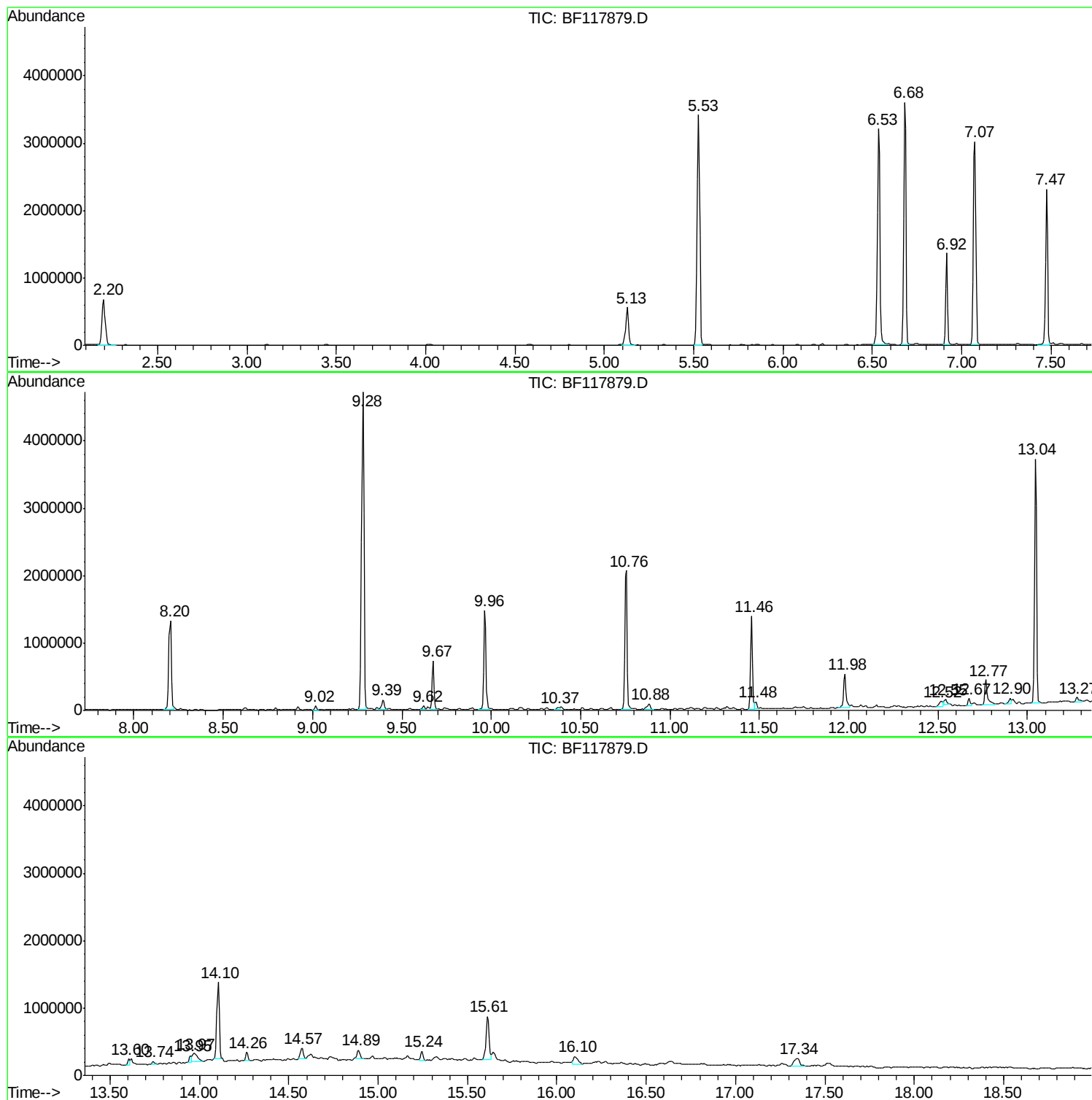
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF111319.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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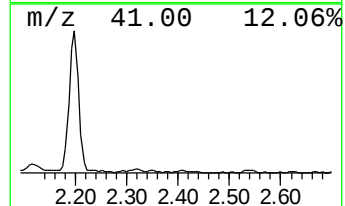
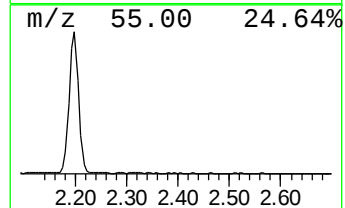
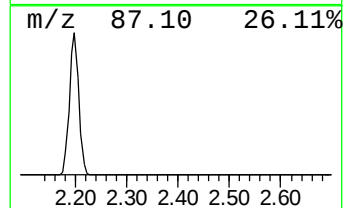
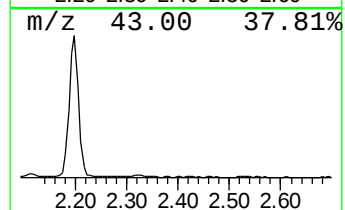
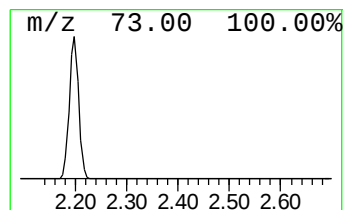
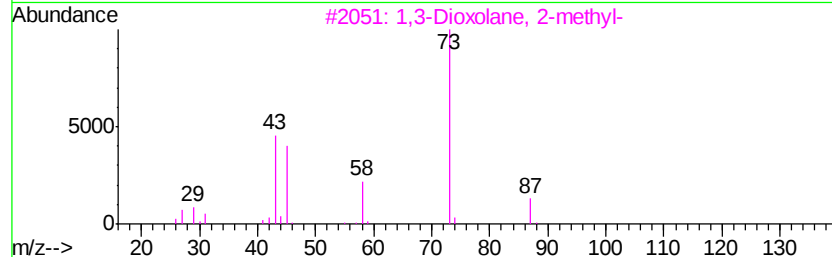
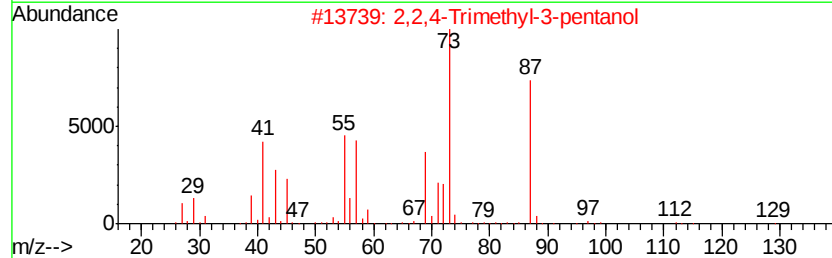
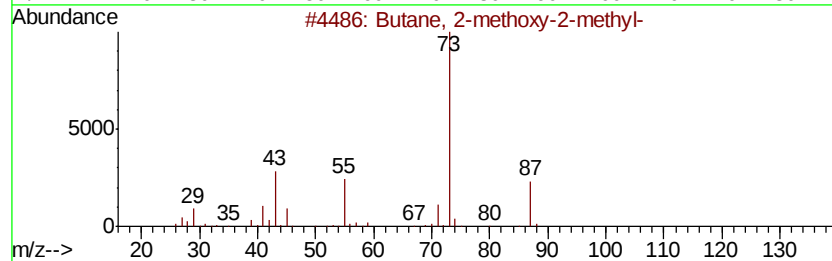
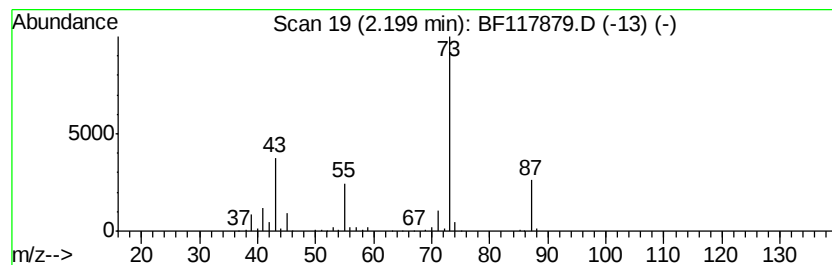
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF111319.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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.
2.20	16.29 ng	854650	1,4-Dichlorobenzene-d4	6.92

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2		2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	40
3		1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	25
4		Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	22
5		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	17



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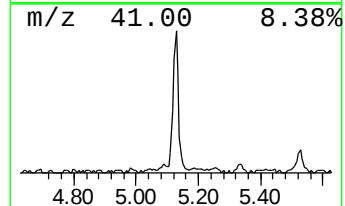
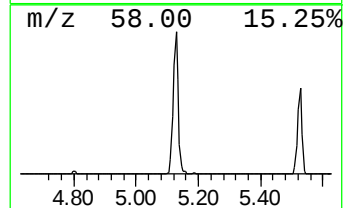
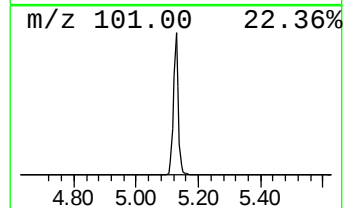
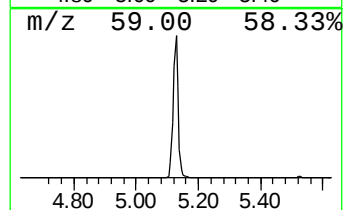
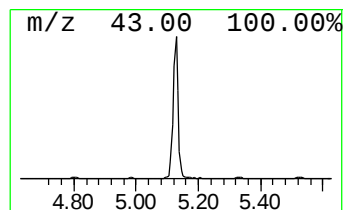
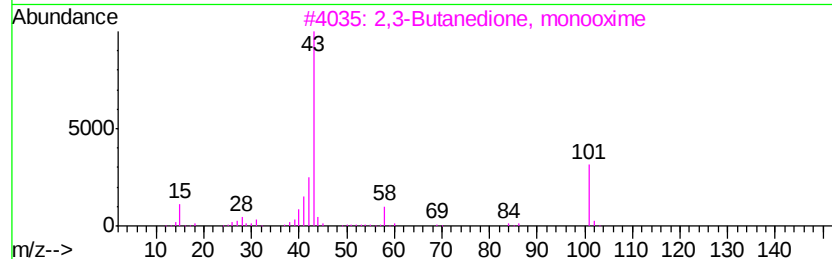
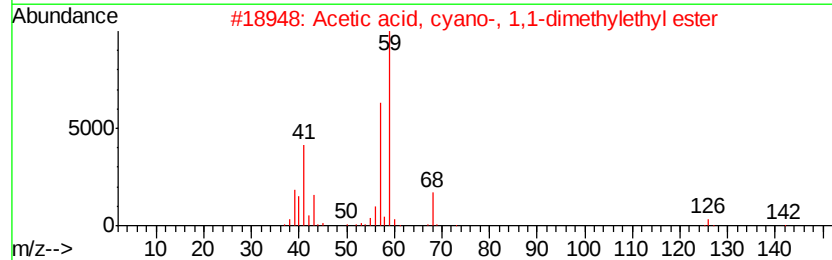
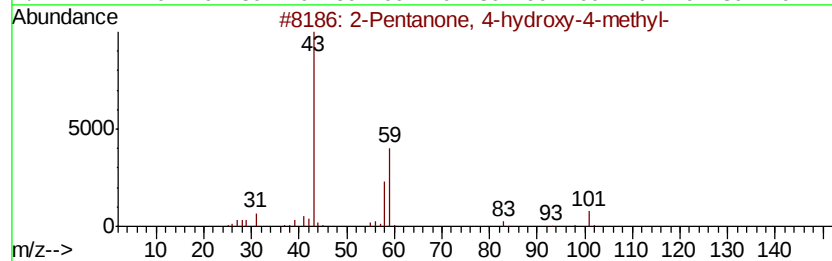
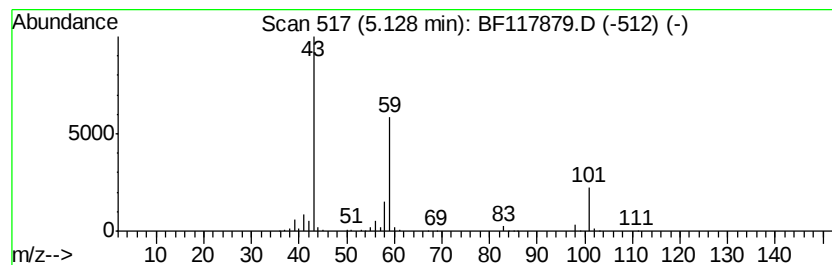
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.13	12.12 ng	635618	1,4-Dichlorobenzene-d4	6.92

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	40
2		Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	32
3		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	25
4		(+)-4-Amino-4,5-dihydro-2(3H)-f...	101	C4H7NO2	016504-58-8	9
5		Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9



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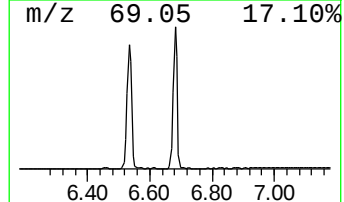
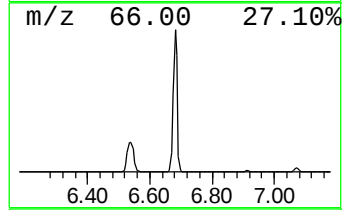
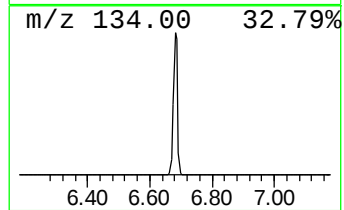
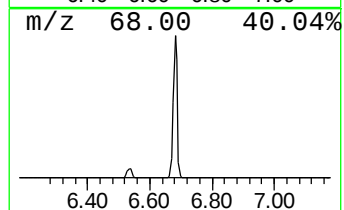
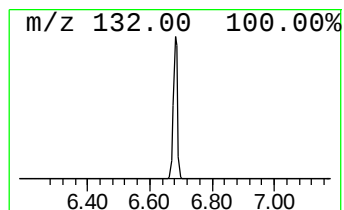
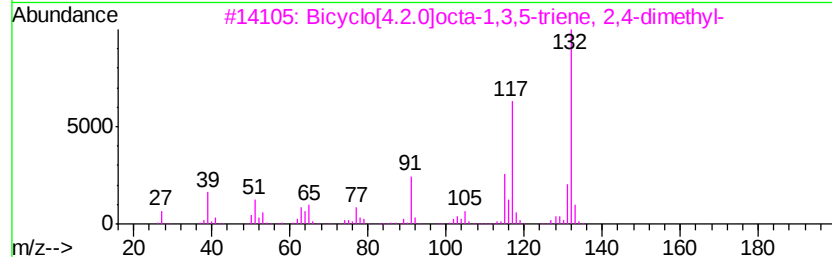
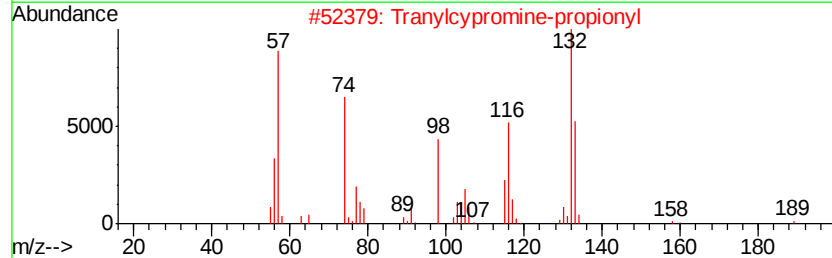
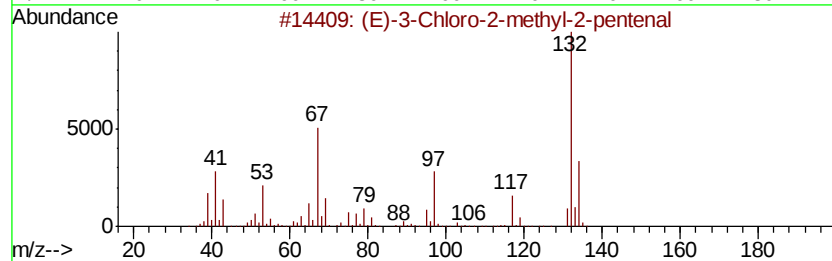
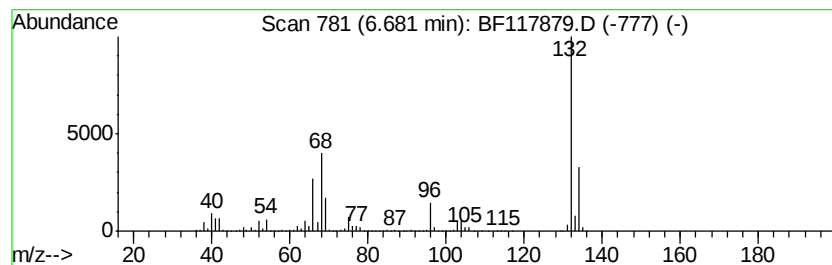
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 3 unknown6.68 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.68	65.46 ng	3434350	1,4-Dichlorobenzene-d4	6.92

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	16
2		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
3		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
4		Benzene, 1-ethenyl-3,5-dimethyl-	132	C10H12	005379-20-4	9
5		Xenon	132	Xe	007440-63-3	9



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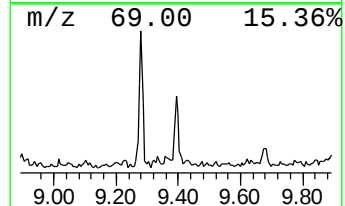
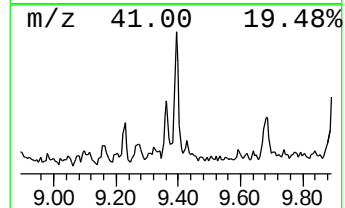
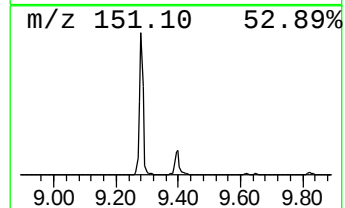
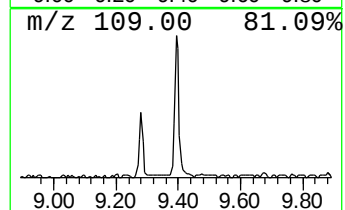
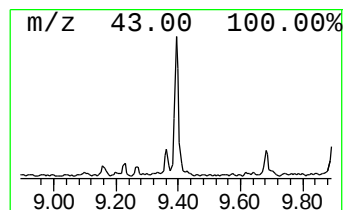
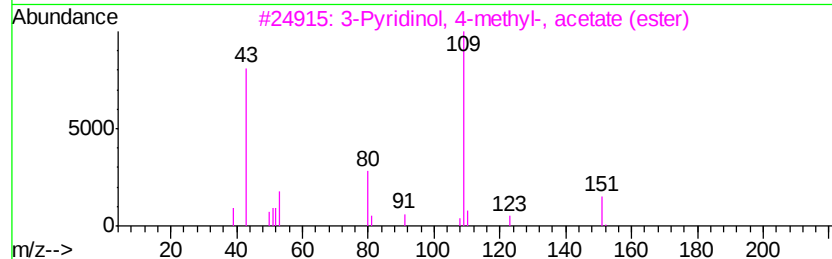
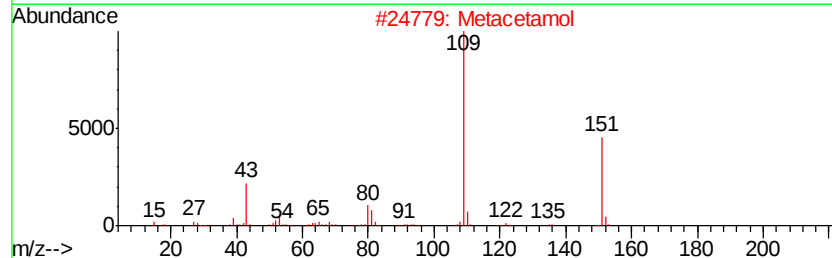
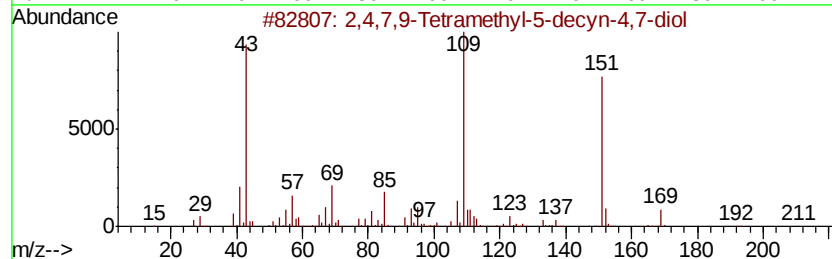
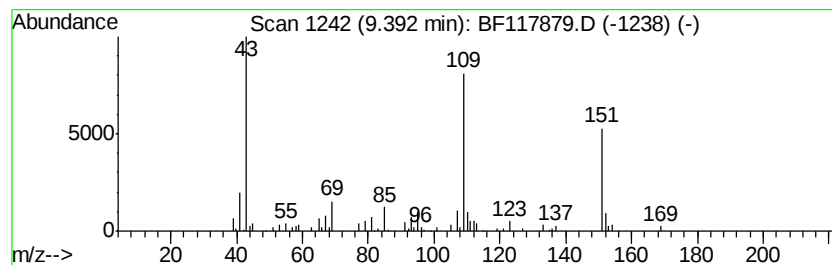
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TIC Library : C:\DATABASE\NIST11.L
 TIC Integration Parameters: LSCINT.P

 Peak Number 5 2,4,7,9-Tetramethyl-5-decyn... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.39	2.09 ng	138225	Acenaphthene-d10	9.96

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2,4,7,9-Tetramethyl-5-decyn-4,7-...	226	C14H26O2	000126-86-3	91
2		Metacetamol	151	C8H9NO2	000621-42-1	59
3		3-Pyridinol, 4-methyl-, acetate ...	151	C8H9NO2	001006-96-8	59
4		p-Isopropoxyvaniline	151	C9H13NO	007664-66-6	59
5		m-Isopropoxyvaniline	151	C9H13NO	041406-00-2	53



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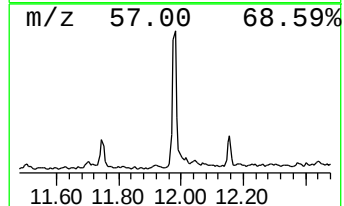
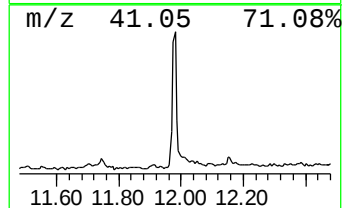
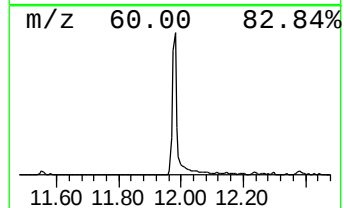
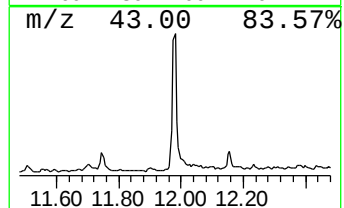
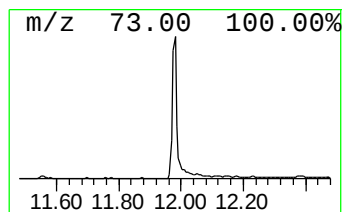
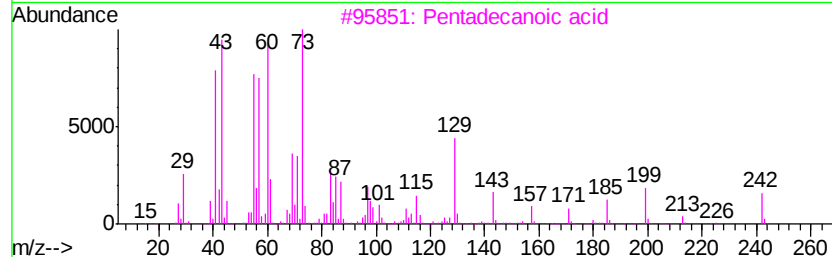
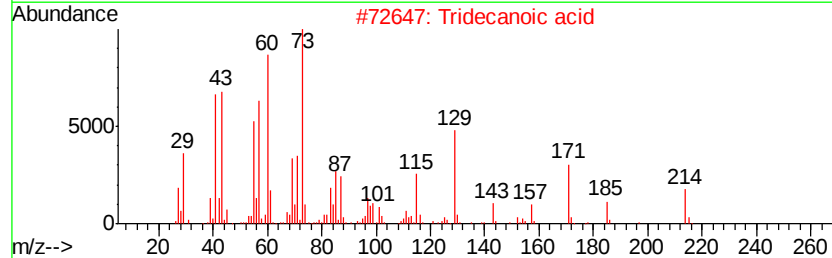
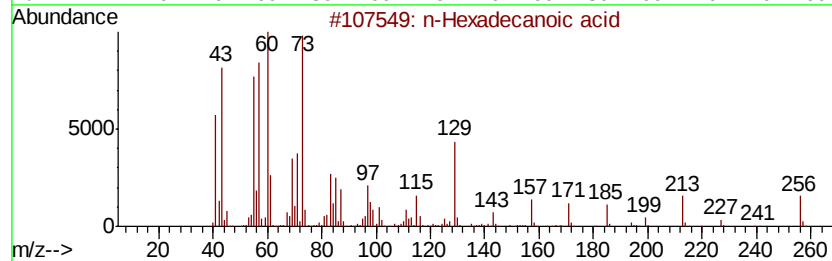
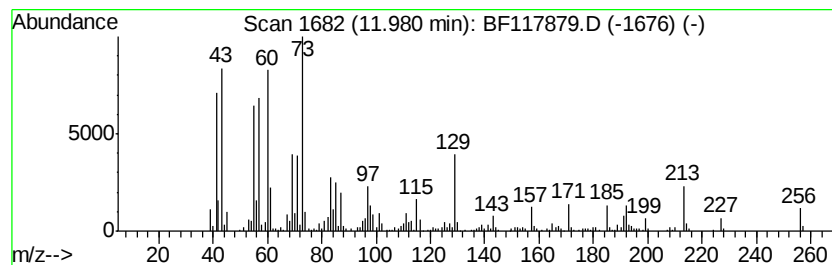
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 n-Hexadecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.98	9.53 ng	552908	Phenanthrene-d10	11.46

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2		Tridecanoic acid	214	C13H26O2	000638-53-9	95
3		Pentadecanoic acid	242	C15H30O2	001002-84-2	93
4		Tetradecanoic acid	228	C14H28O2	000544-63-8	80
5		n-Decanoic acid	172	C10H20O2	000334-48-5	76



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF111919\
Data File : BF117879.D
Acq On : 20 Nov 2019 3:27
Operator : JU
Sample : K5908-02
Misc :
ALS Vial : 27 Sample Multiplier: 1

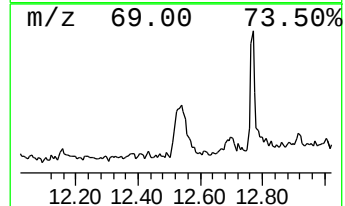
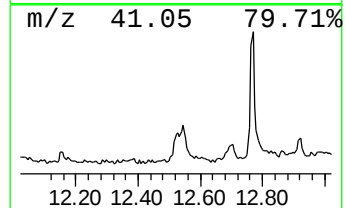
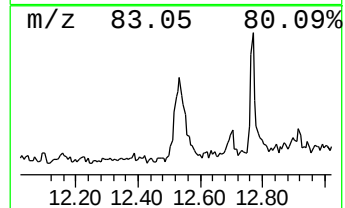
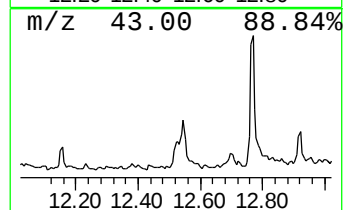
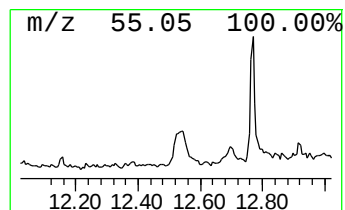
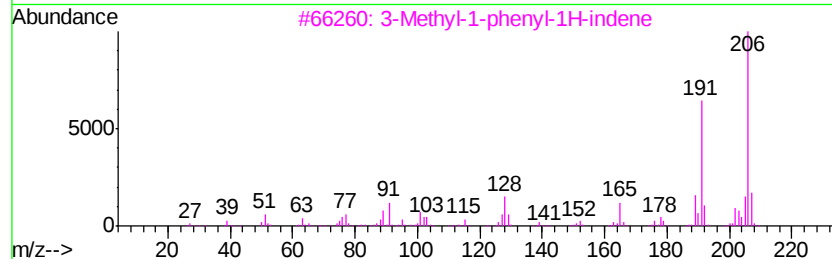
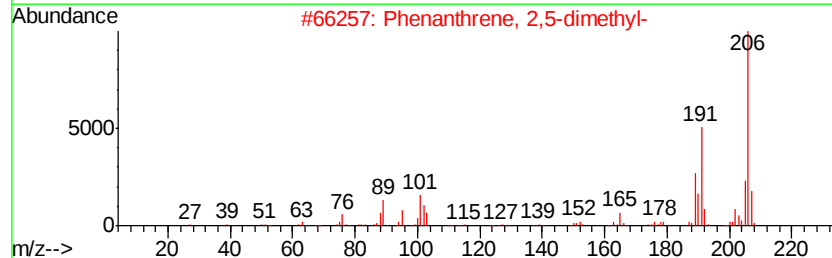
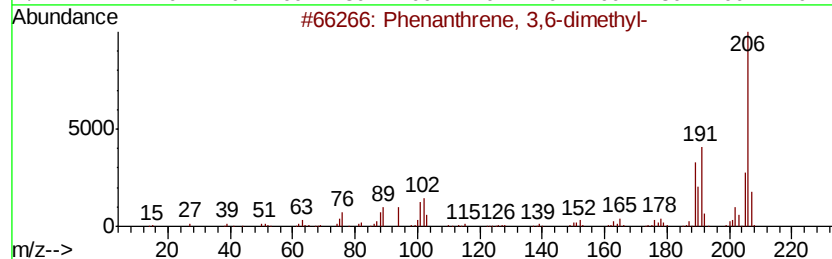
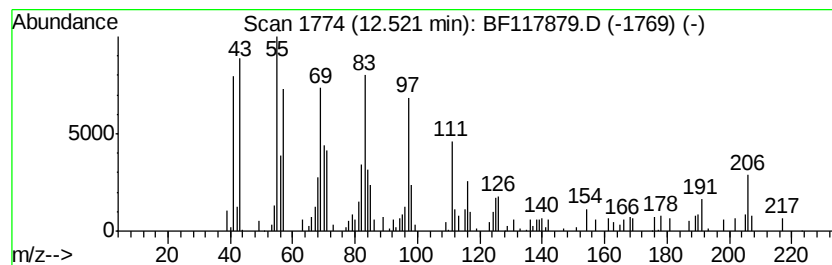
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ClientSampleId :
B-8-1-(6-12)

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF111319.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Phenanthrene, 3,6-dimethyl- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.52	2.00 ng	116079	Phenanthrene-d10	11.46	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 3,6-dimethyl-	206	C16H14	001576-67-6	70
2	Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	58
3	3-Methyl-1-phenyl-1H-indene	206	C16H14	022360-63-0	55
4	Naphthalene, 1,2-dihydro-4-phenyl-	206	C16H14	007469-40-1	49
5	Phenanthrene, 1,7-dimethyl-	206	C16H14	000483-87-4	46



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF111919\
Data File : BF117879.D
Acq On : 20 Nov 2019 3:27
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Misc :
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Instrument :
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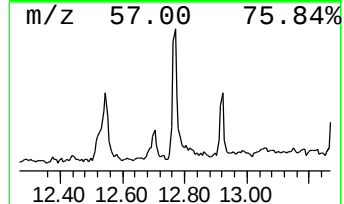
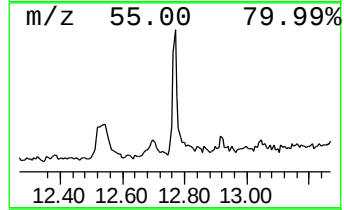
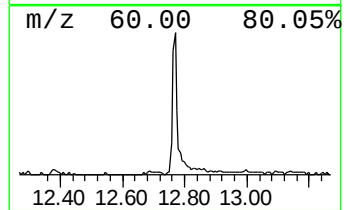
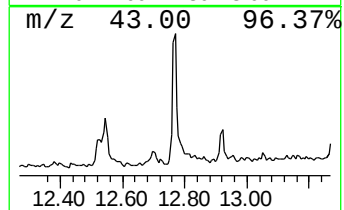
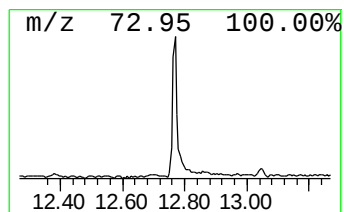
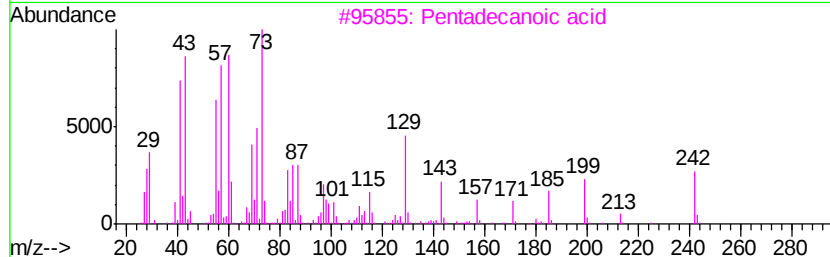
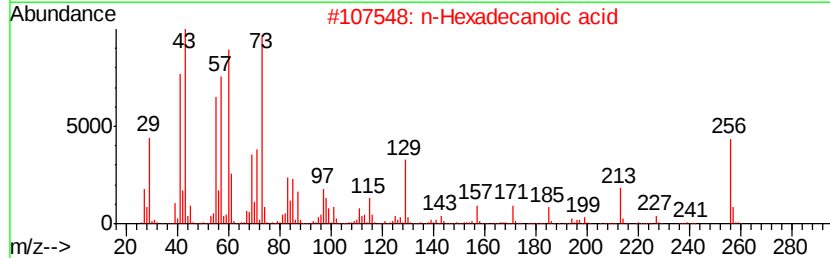
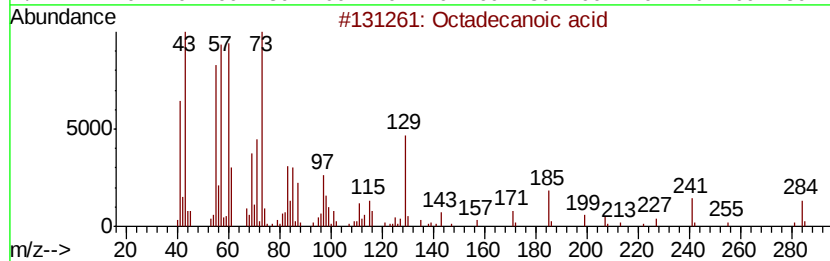
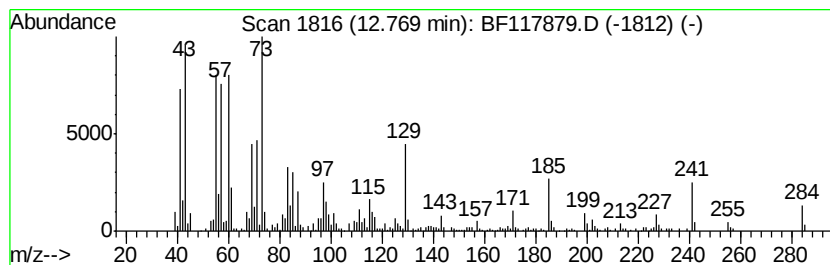
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Octadecanoic acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.77	7.96 ng	461338	Phenanthrene-d10	11.46

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Octadecanoic acid	284	C18H36O2	000057-11-4	99
2		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	93
3		Pentadecanoic acid	242	C15H30O2	001002-84-2	89
4		Tetradecanoic acid	228	C14H28O2	000544-63-8	83
5		Tridecanoic acid	214	C13H26O2	000638-53-9	76



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF111919\
Data File : BF117879.D
Acq On : 20 Nov 2019 3:27
Operator : JU
Sample : K5908-02
Misc :
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ClientSampleId :
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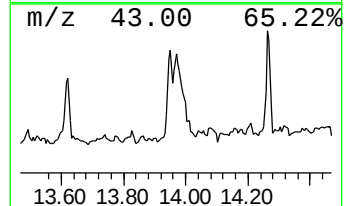
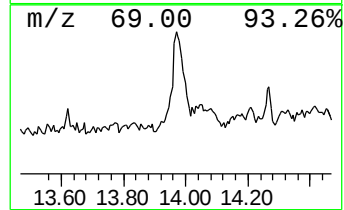
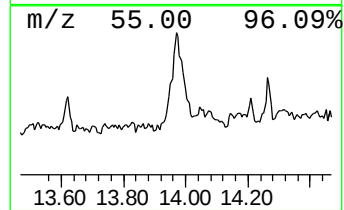
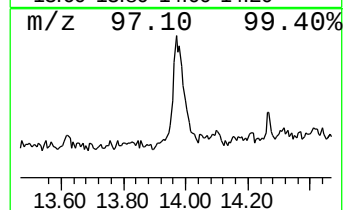
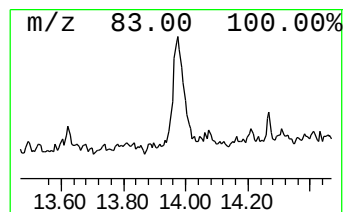
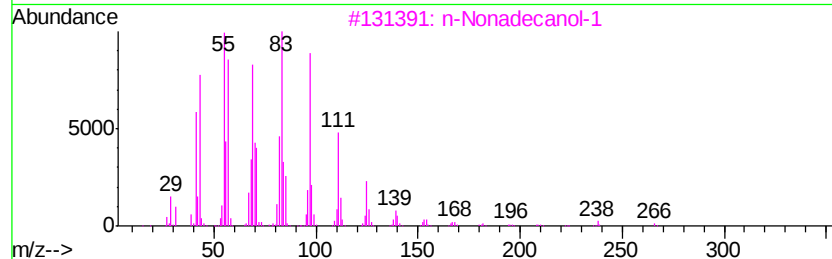
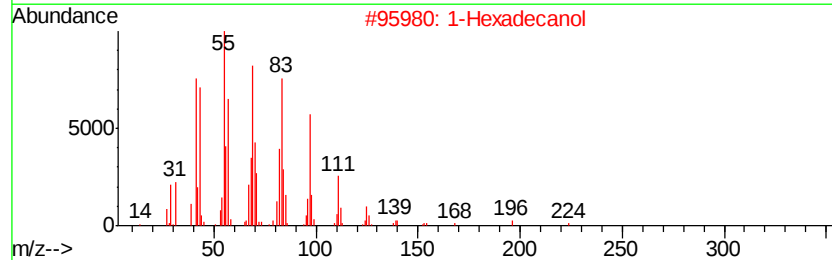
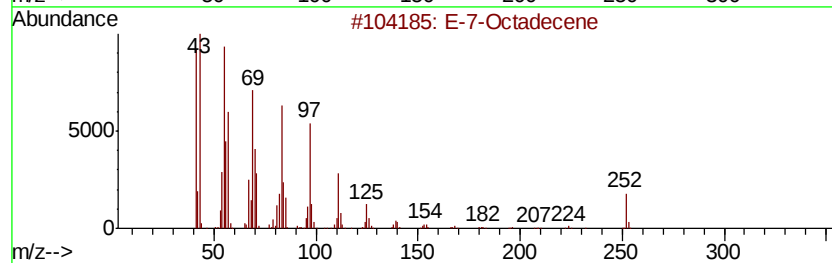
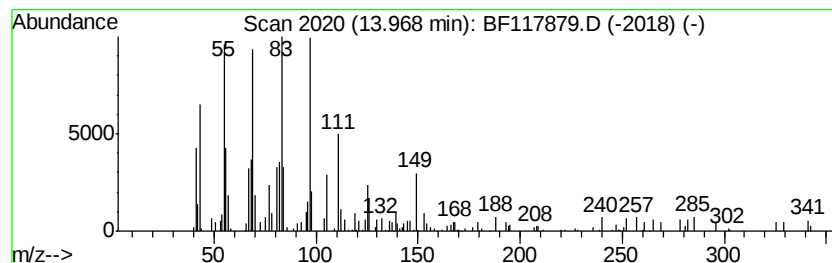
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 E-7-Octadecene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.97	5.15 ng	252542	Chrysene-d12	14.10

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	E-7-Octadecene	252	C18H36	1000130-92-0	89
2		1-Hexadecanol	242	C16H34O	036653-82-4	64
3		n-Nonadecanol-1	284	C19H40O	001454-84-8	58
4		Decane, 5,6-bis(2,2-dimethylprop...	278	C20H38	073002-85-4	58
5		1-Heneicosanol	312	C21H44O	015594-90-8	52



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF111919\
Data File : BF117879.D
Acq On : 20 Nov 2019 3:27
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Misc :
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Instrument :
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ClientSampleId :
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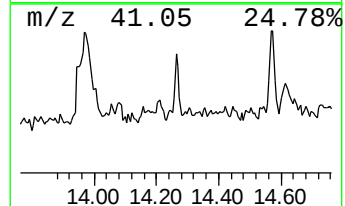
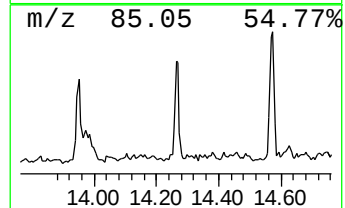
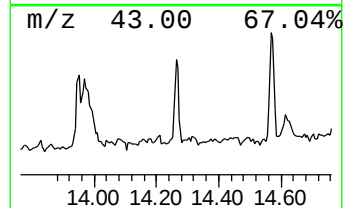
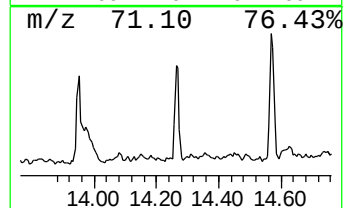
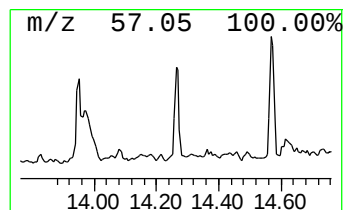
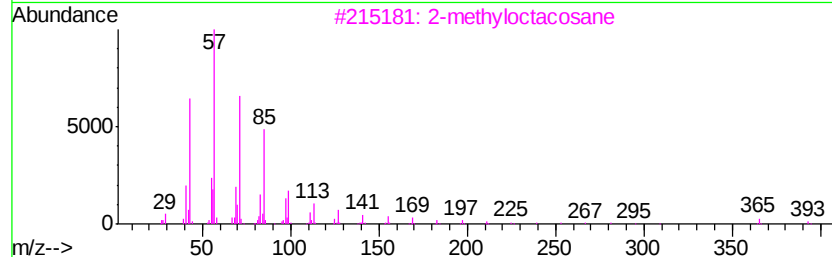
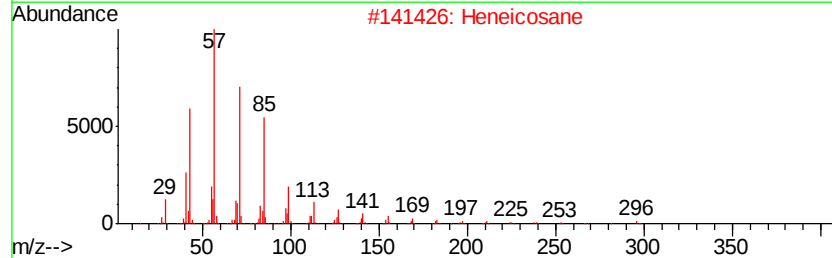
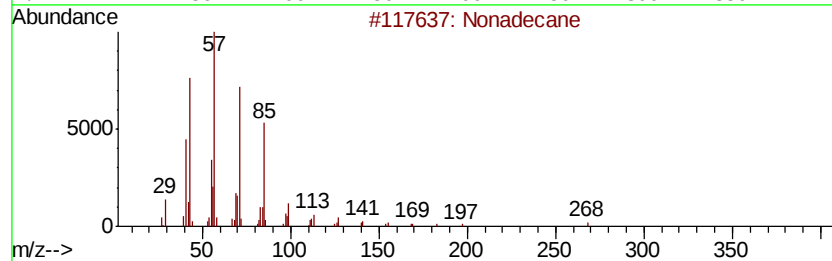
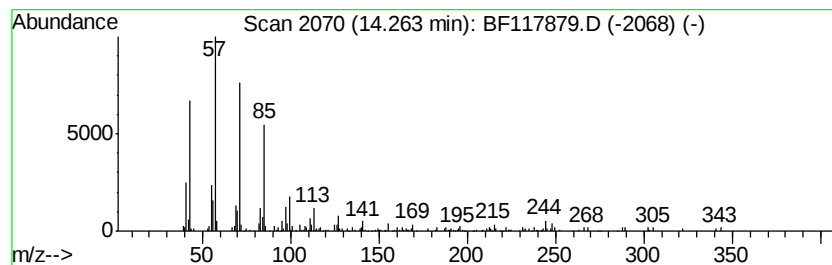
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Nonadecane Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.26	2.47 ng	121220	Chrysene-d12	14.10

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Nonadecane	268	C19H40	000629-92-5	91
2		Heneicosane	296	C21H44	000629-94-7	90
3		2-methyloctacosane	408	C29H60	1000376-72-8	90
4		Hexacosane	366	C26H54	000630-01-3	90
5		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF111919\
Data File : BF117879.D
Acq On : 20 Nov 2019 3:27
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Sample : K5908-02
Misc :
ALS Vial : 27 Sample Multiplier: 1

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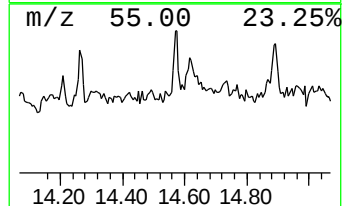
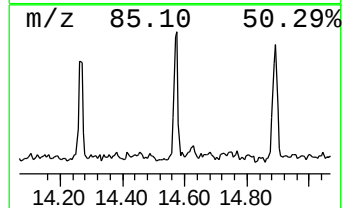
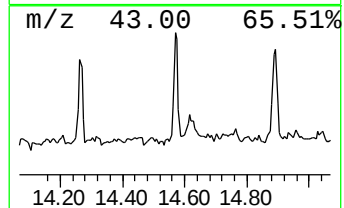
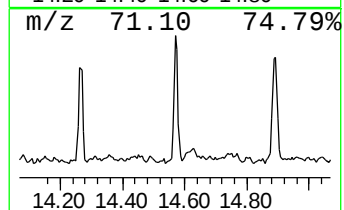
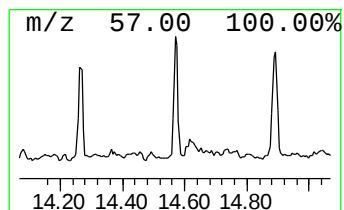
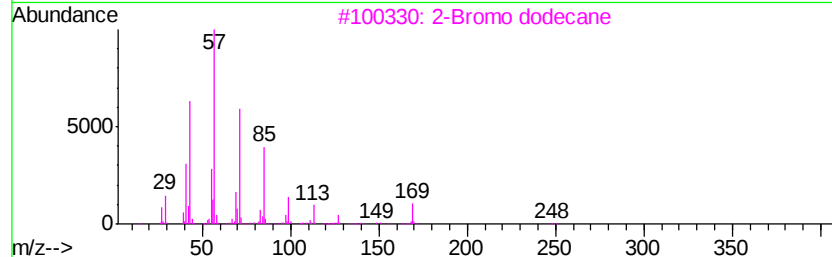
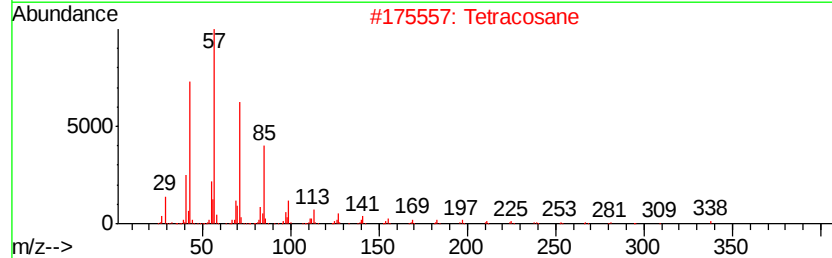
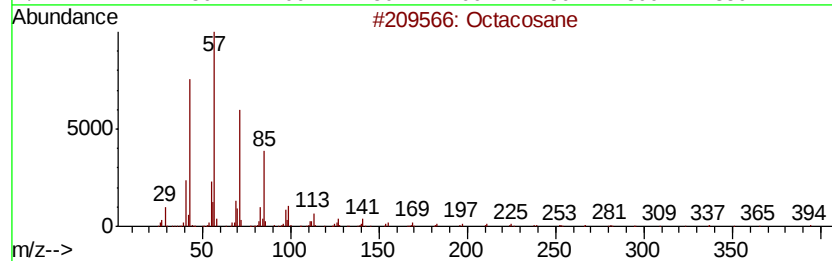
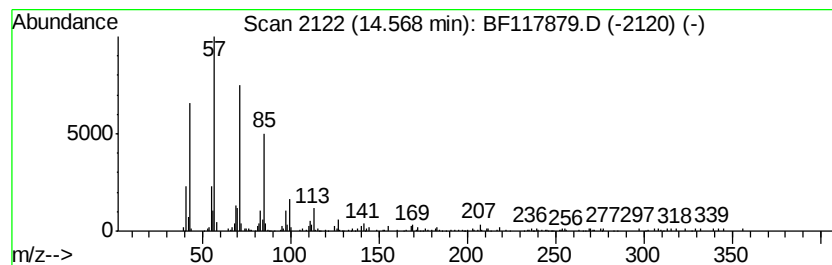
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Octacosane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.57	3.03 ng	148437	Chrysene-d12	14.10

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octacosane	394	C28H58	000630-02-4	91
2		Tetracosane	338	C24H50	000646-31-1	91
3		2-Bromo dodecane	248	C12H25Br	013187-99-0	90
4		Pentacosane	352	C25H52	000629-99-2	90
5		Octadecane, 1-iodo-	380	C18H37I	000629-93-6	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF111919\
Data File : BF117879.D
Acq On : 20 Nov 2019 3:27
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Misc :
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ClientSampleId :
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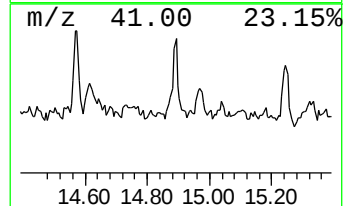
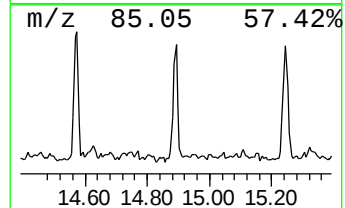
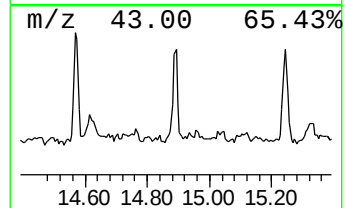
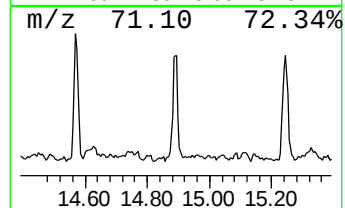
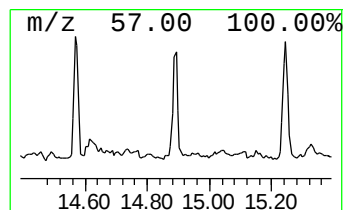
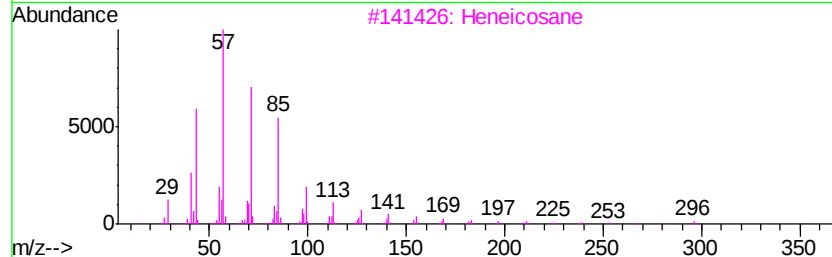
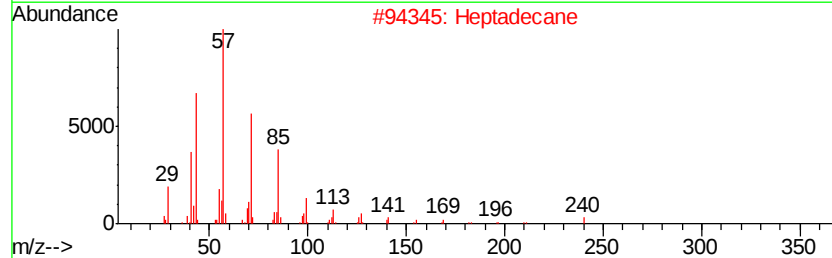
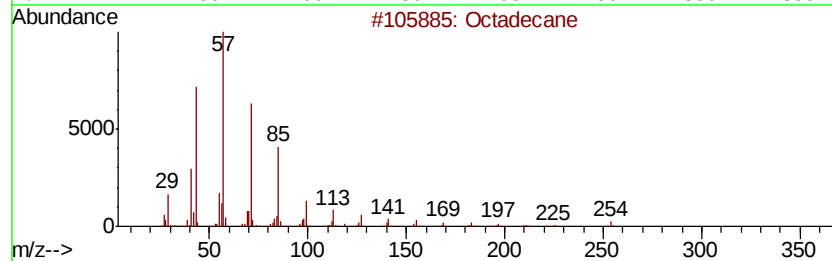
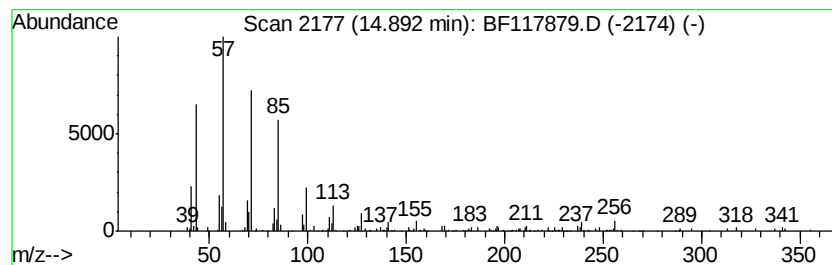
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Octadecane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.89	3.17 ng	127342	Perylene-d12	15.61

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecane	254	C18H38	000593-45-3	93
2		Heptadecane	240	C17H36	000629-78-7	87
3		Heneicosane	296	C21H44	000629-94-7	87
4		Tridecane, 7-hexyl-	268	C19H40	007225-66-3	87
5		Tridecanol, 2-ethyl-2-methyl-	242	C16H34O	1000115-66-1	86



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF111919\
Data File : BF117879.D
Acq On : 20 Nov 2019 3:27
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Misc :
ALS Vial : 27 Sample Multiplier: 1

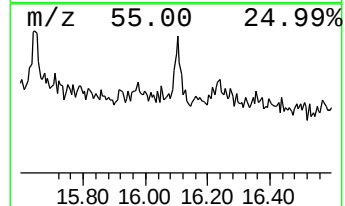
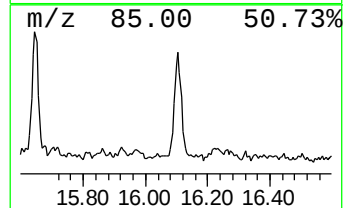
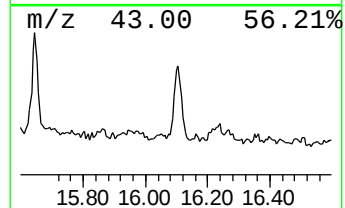
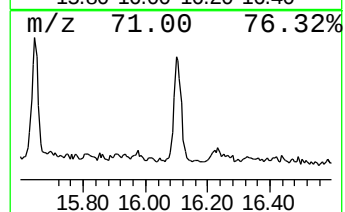
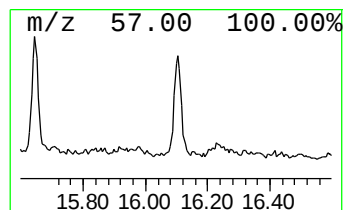
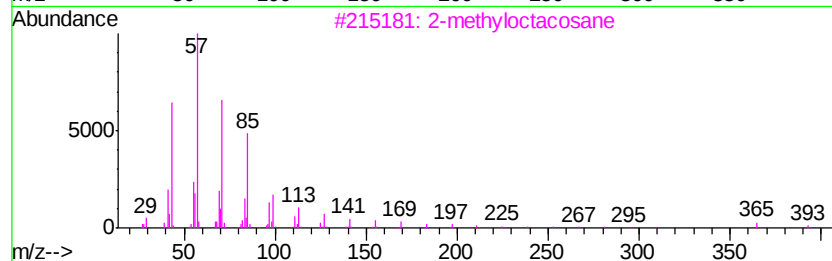
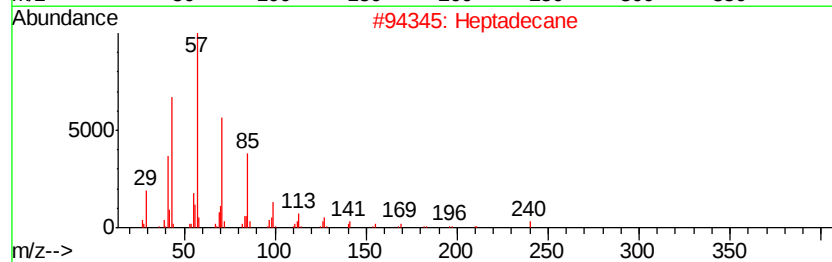
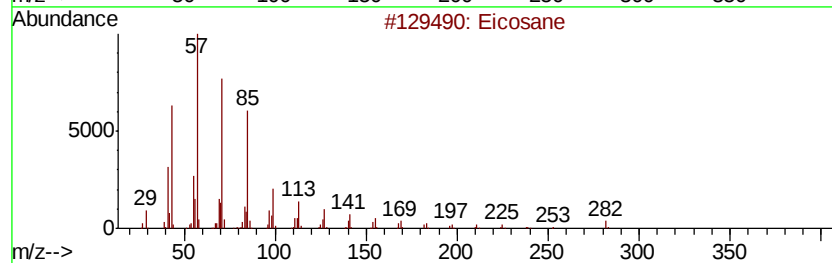
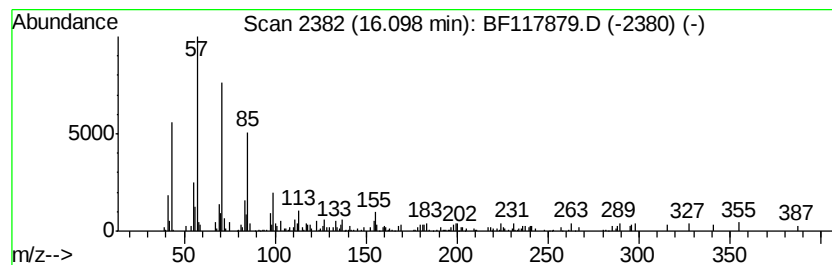
Instrument :
BNA_F
ClientSampleId :
B-8-1-(6-12)

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF111319.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Eicosane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.10	4.76 ng	190909	Perylene-d12	15.61
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Eicosane	282 C20H42	000112-95-8	90
2	Heptadecane	240 C17H36	000629-78-7	90
3	2-methyloctacosane	408 C29H60	1000376-72-8	81
4	Hexadecane, 2-methyl-	240 C17H36	001560-92-5	74
5	Heptacosane	380 C27H56	000593-49-7	74



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF111919\
Data File : BF117879.D
Acq On : 20 Nov 2019 3:27
Operator : JU
Sample : K5908-02
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
B-8-1-(6-12)

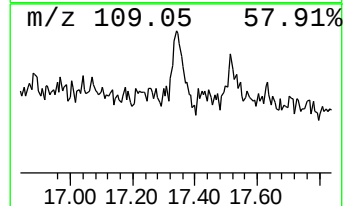
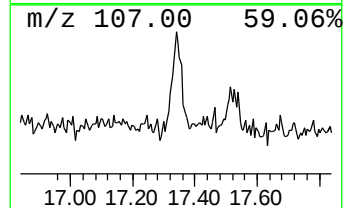
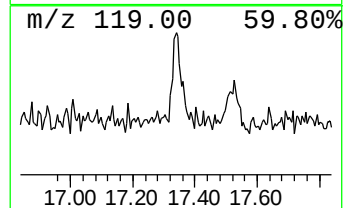
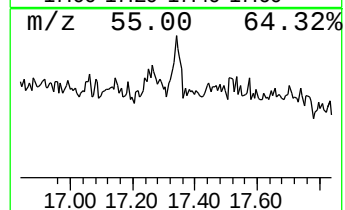
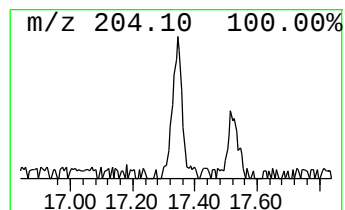
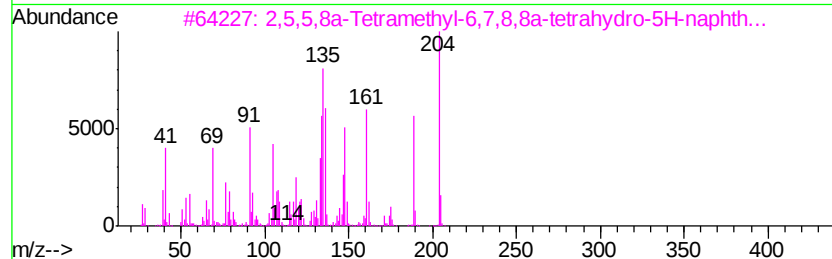
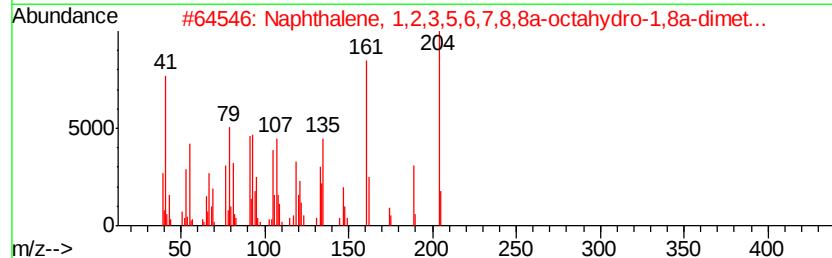
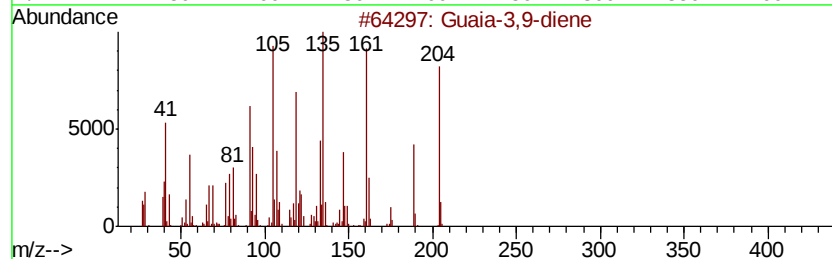
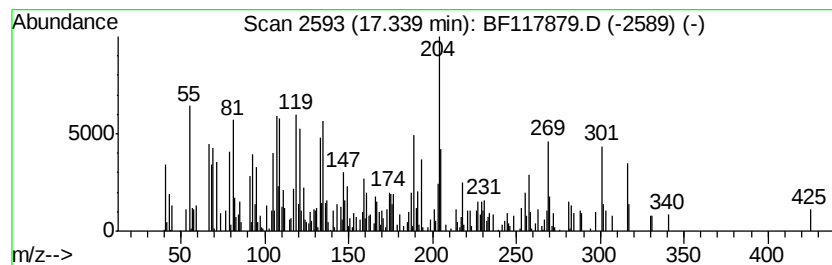
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF111319.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 Guaia-3,9-diene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.34	6.10 ng	245060	Perylene-d12	15.61

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Guaia-3,9-diene	204	C15H24	000489-83-8	62
2			Naphthalene, 1,2,3,5,6,7,8,8a-oc...	204	C15H24	004630-07-3	58
3			2,5,5,8a-Tetramethyl-6,7,8,8a-te...	204	C14H200	124957-09-1	50
4			1H-3a,7-Methanoazulene, octahydr...	204	C15H24	000508-55-4	49
5			1,4-Methanoazulen-9-ol, decahydr...	222	C15H260	000465-24-7	46



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF111919\
 Data File : BF117879.D
 Acq On : 20 Nov 2019 3:27
 Operator : JU
 Sample : K5908-02
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 B-8-1-(6-12)

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF111319.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...	2.20	16.3	ng	854650	1	6.92	1049240	20.0
2-Pentanone, 4-hy...	5.13	12.1	ng	635618	1	6.92	1049240	20.0
unknown6.68	6.68	65.5	ng	3434350	1	6.92	1049240	20.0
2,4,7,9-Tetrameth...	9.39	2.1	ng	138225	3	9.96	1320000	20.0
n-Hexadecanoic acid	11.98	9.5	ng	552908	4	11.46	1159840	20.0
Phenanthrene, 3,6...	12.52	2.0	ng	116079	4	11.46	1159840	20.0
Octadecanoic acid	12.77	8.0	ng	461338	4	11.46	1159840	20.0
E-7-Octadecene	13.97	5.2	ng	252542	5	14.10	980623	20.0
Nonadecane	14.26	2.5	ng	121220	5	14.10	980623	20.0
Octacosane	14.57	3.0	ng	148437	5	14.10	980623	20.0
Octadecane	14.89	3.2	ng	127342	6	15.61	802915	20.0
Eicosane	16.10	4.8	ng	190909	6	15.61	802915	20.0
Guaia-3,9-diene	17.34	6.1	ng	245060	6	15.61	802915	20.0