

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF042219\
 Data File : BF113892.D
 Acq On : 22 Apr 2019 21:50
 Operator : JU/SJ
 Sample : K2515-19
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-13

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-BF042219.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	2.216	19	22	35	rVB2	62588	108516	2.05%	0.265%
2	5.522	579	584	587	rBV	3920489	4293515	81.19%	10.479%
3	6.528	748	755	757	rBV	3669417	4366396	82.57%	10.657%
4	6.669	774	779	782	rBV	4720509	4543976	85.93%	11.091%
5	6.892	814	817	821	rBV	811390	679208	12.84%	1.658%
6	7.051	840	844	848	rBV	3894361	3743839	70.80%	9.138%
7	7.451	907	912	915	rBV	2763284	2813231	53.20%	6.866%
8	8.175	1031	1035	1038	rBV	995459	788301	14.91%	1.924%
9	9.251	1213	1218	1221	rBV	5848670	5288066	100.00%	12.907%
10	9.569	1269	1272	1275	rBV	176283	152147	2.88%	0.371%
11	9.639	1280	1284	1287	rBV	348458	285849	5.41%	0.698%
12	9.928	1329	1333	1337	rBV	1067339	861851	16.30%	2.104%
13	10.716	1462	1467	1470	rBV	3343225	2966782	56.10%	7.241%
14	11.410	1581	1585	1588	rBV	947367	787192	14.89%	1.921%
15	11.933	1671	1674	1686	rVB2	336008	357708	6.76%	0.873%
16	12.622	1787	1791	1794	rBV	181411	155648	2.94%	0.380%
17	12.722	1804	1808	1815	rBV	175593	184150	3.48%	0.449%
18	12.851	1825	1830	1832	rBV	139713	133801	2.53%	0.327%
19	12.998	1850	1855	1858	rBV	4933677	4501960	85.13%	10.988%
20	13.433	1925	1929	1932	rBV	185331	221586	4.19%	0.541%
21	13.463	1932	1934	1942	rVB3	68537	117732	2.23%	0.287%
22	13.974	2014	2021	2028	rBV6	59590	184096	3.48%	0.449%
23	14.045	2028	2033	2036	rVV	952857	957965	18.12%	2.338%
24	14.068	2036	2037	2043	rVB	98632	72881	1.38%	0.178%
25	14.221	2059	2063	2066	rVB3	82715	85723	1.62%	0.209%
26	14.527	2112	2115	2118	rVB2	64350	59783	1.13%	0.146%
27	14.833	2164	2167	2171	rVB	75588	74426	1.41%	0.182%
28	14.915	2177	2181	2187	rBV2	50021	74234	1.40%	0.181%
29	15.092	2207	2211	2221	rVB	113032	216993	4.10%	0.530%
30	15.180	2222	2226	2232	rVB2	103950	133853	2.53%	0.327%
31	15.398	2259	2263	2269	rVB	69519	91255	1.73%	0.223%
32	15.457	2269	2273	2278	rVV	91356	118489	2.24%	0.289%
33	15.521	2279	2284	2288	rVV	758712	981208	18.56%	2.395%
34	15.568	2288	2292	2297	rVB2	92446	141780	2.68%	0.346%

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ALS Vial : 10 Sample Multiplier: 1

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

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

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35	16.010	2363	2367	2376	rBV2	92856	149134	2.82%	0.364%
36	16.527	2450	2455	2459	rBV2	54859	103364	1.95%	0.252%
37	16.998	2530	2535	2543	rVB2	51098	104971	1.99%	0.256%
38	17.445	2606	2611	2618	rVB	38437	70026	1.32%	0.171%

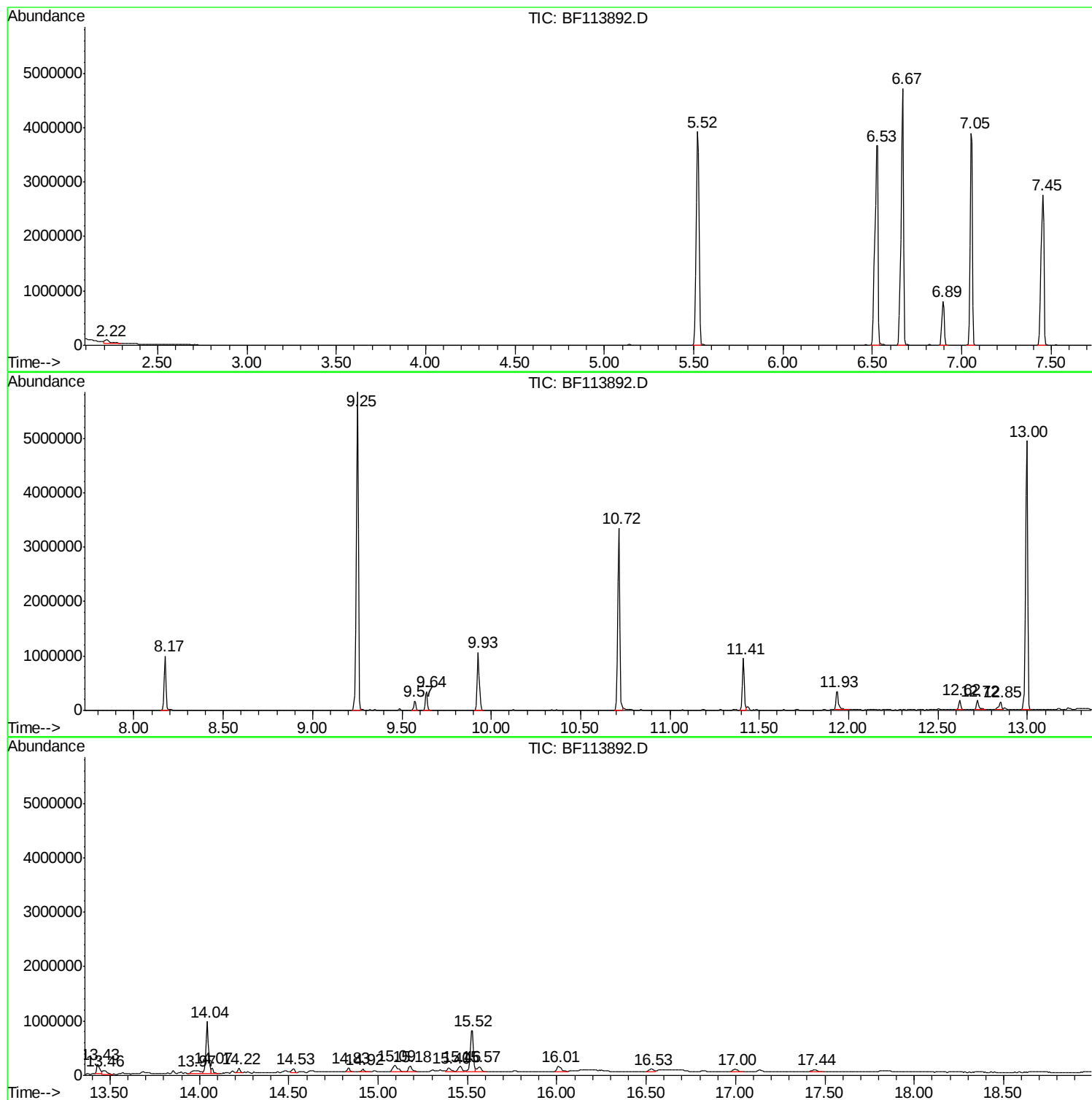
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF042219.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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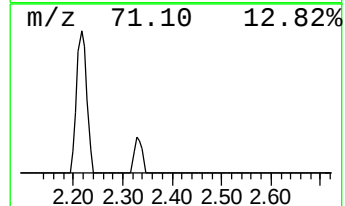
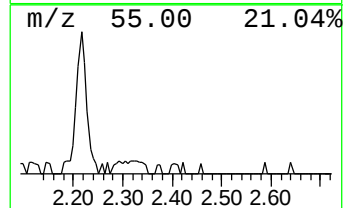
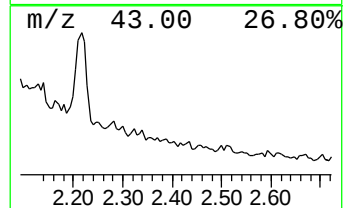
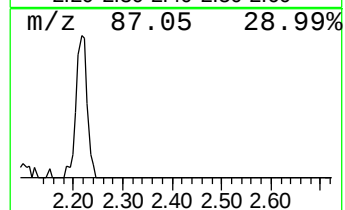
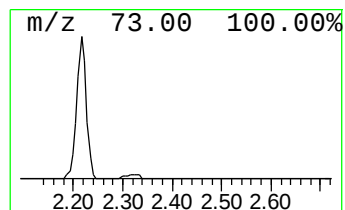
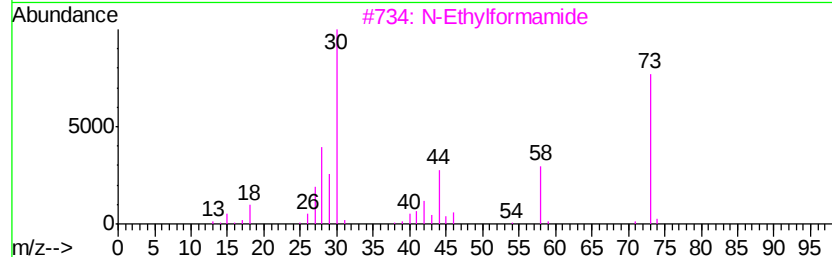
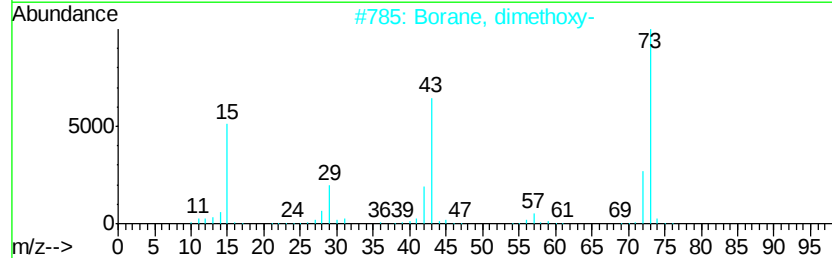
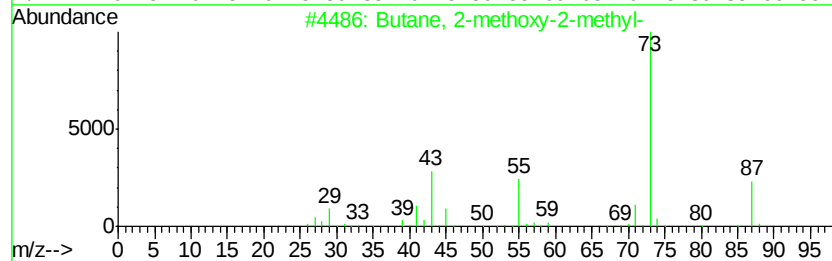
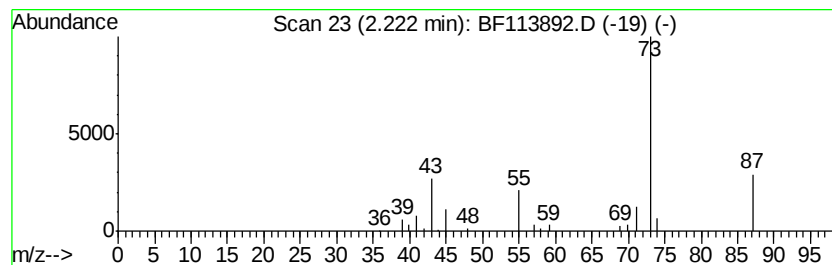
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.	
2.22	3.20 ng	108516	1,4-Dichlorobenzene-d4	6.89	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2	Borane, dimethoxy-	74	C2H7BO2	004542-61-4	9
3	N-Ethylformamide	73	C3H7NO	000627-45-2	9
4	3-Pentanol, 2,4-dimethyl-	116	C7H16O	000600-36-2	9
5	Propane, 2-methoxy-2-methyl-	88	C5H12O	001634-04-4	9



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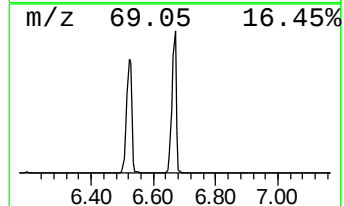
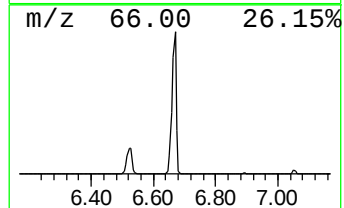
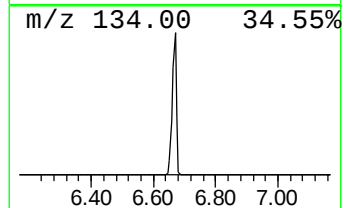
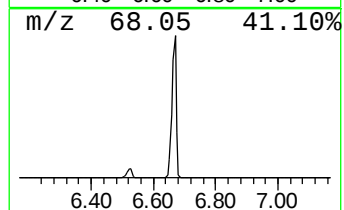
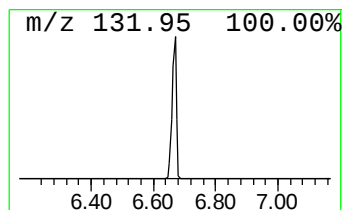
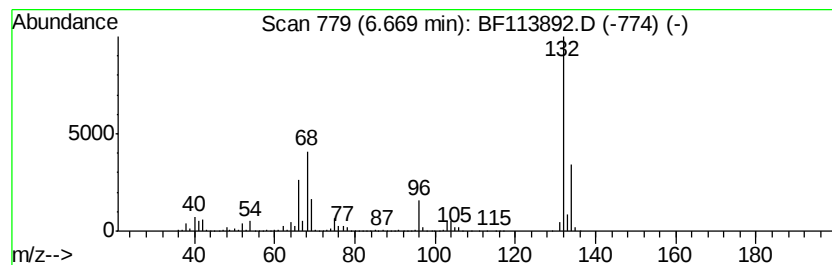
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF042219.M
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 unknown6.67 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.67	133.80 ng	4543980	1,4-Dichlorobenzene-d4	6.89

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	16
3		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	12
4		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
5		2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9



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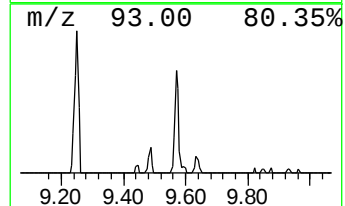
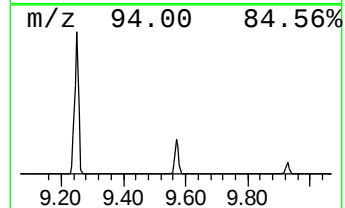
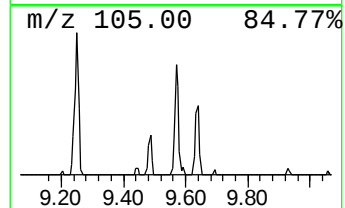
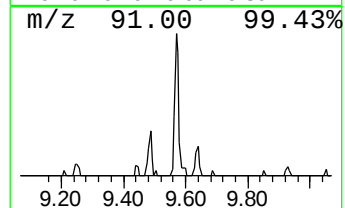
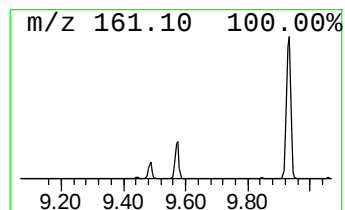
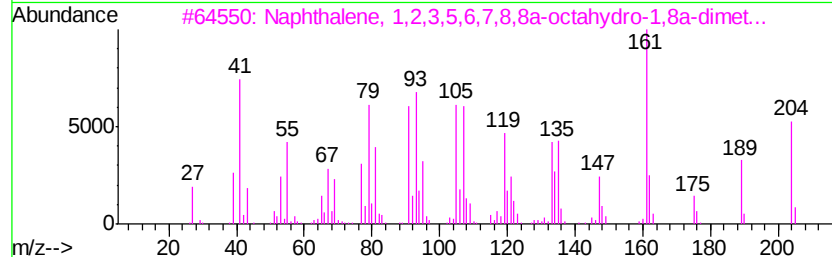
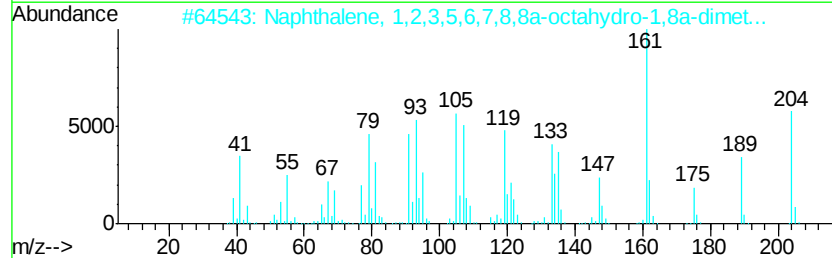
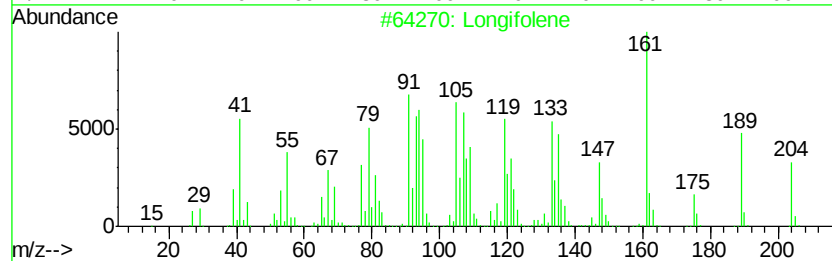
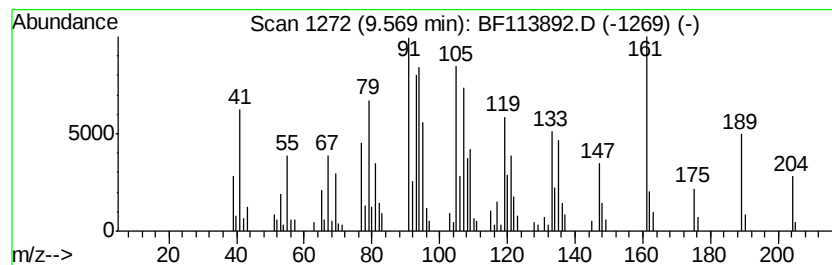
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Longifolene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.57	3.53 ng	152147	Acenaphthene-d10	9.93

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Longifolene	204	C15H24	000475-20-7	99
2		Naphthalene, 1,2,3,5,6,7,8,8a-oc...	204	C15H24	004630-07-3	95
3		Naphthalene, 1,2,3,5,6,7,8,8a-oc...	204	C15H24	010219-75-7	92
4		1H-Cycloprop[e]azulene, decahydr...	204	C15H24	072747-25-2	89
5		Aromandendrene	204	C15H24	000489-39-4	89



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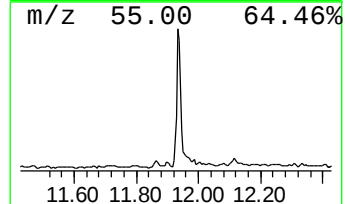
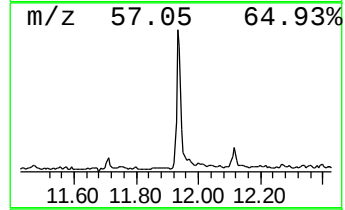
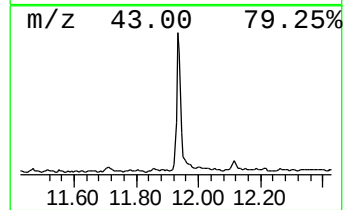
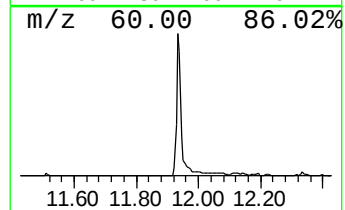
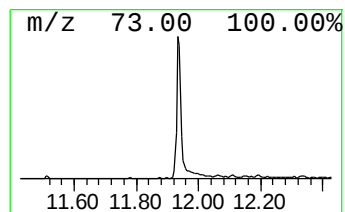
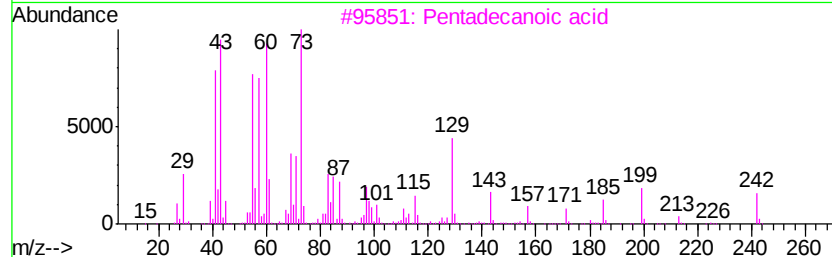
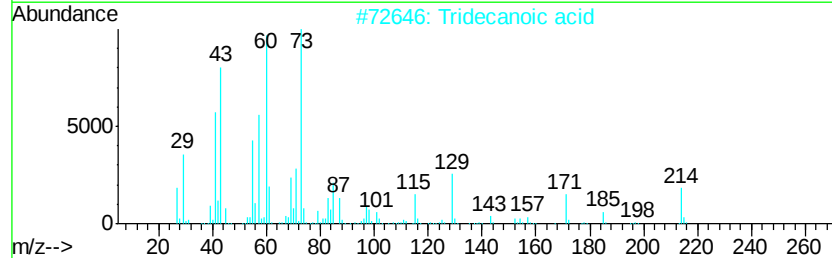
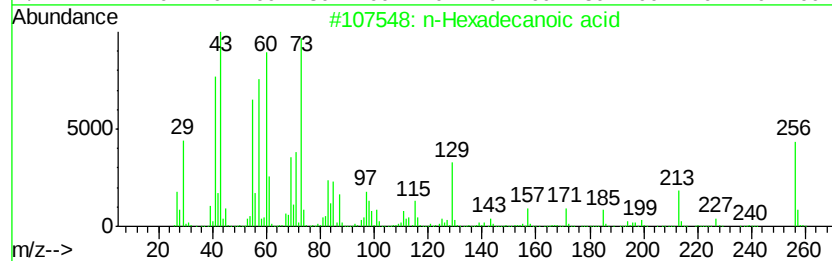
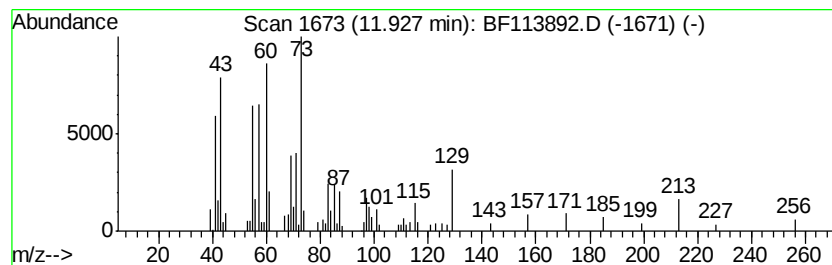
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.93	9.09 ng	357708	Phenanthrene-d10	11.41

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
2		Tridecanoic acid	214	C13H26O2	000638-53-9	91
3		Pentadecanoic acid	242	C15H30O2	001002-84-2	90
4		Tetradecanoic acid	228	C14H28O2	000544-63-8	90
5		Oxalic acid, dodecyl propyl ester	300	C17H32O4	1000309-26-5	18



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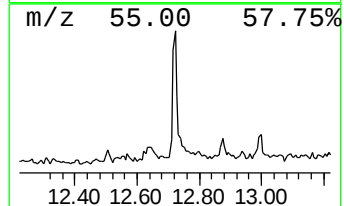
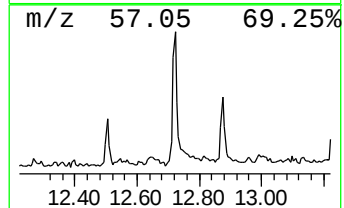
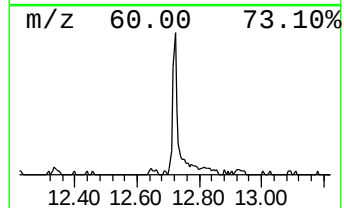
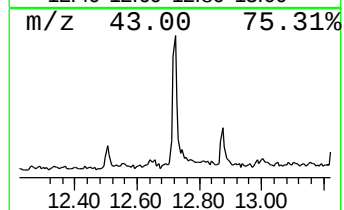
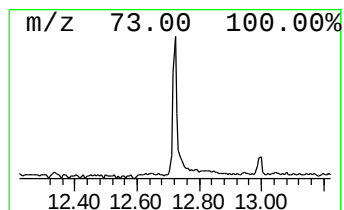
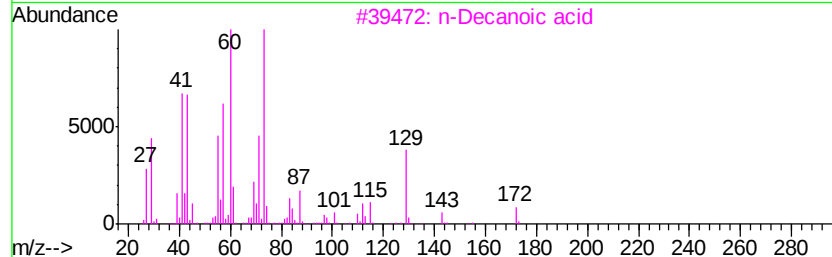
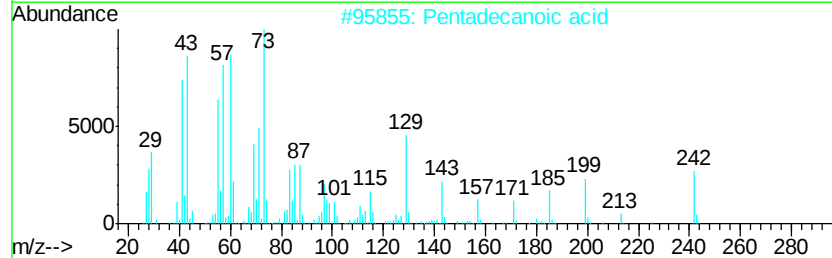
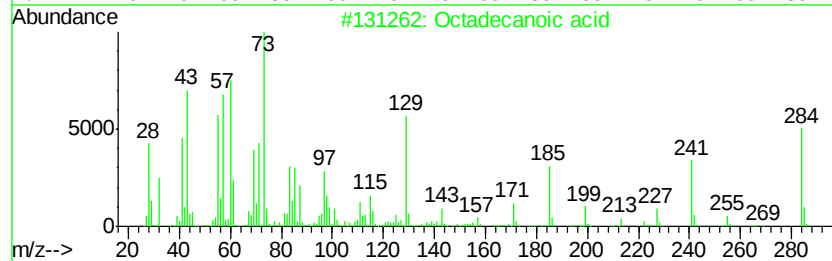
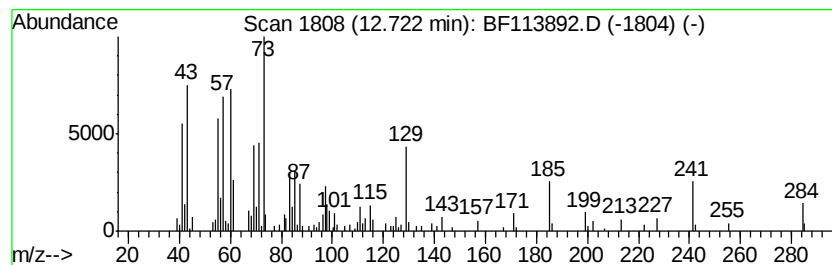
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 6 Octadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.72	4.68 ng	184150	Phenanthrene-d10	11.41

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Octadecanoic acid	284	C18H36O2	000057-11-4	99
2		Pentadecanoic acid	242	C15H30O2	001002-84-2	89
3		n-Decanoic acid	172	C10H20O2	000334-48-5	60
4		Tetradecanoic acid	228	C14H28O2	000544-63-8	58
5		17-Pentatriacontene	491	C35H70	006971-40-0	25



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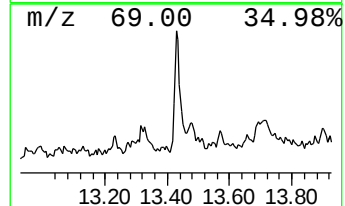
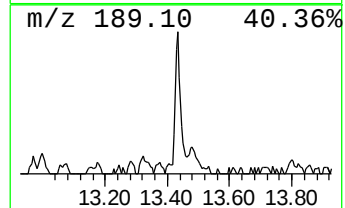
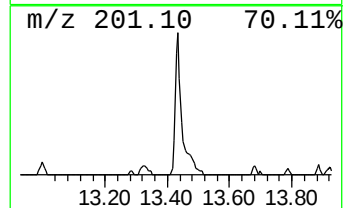
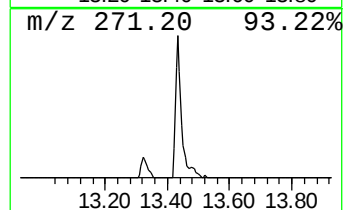
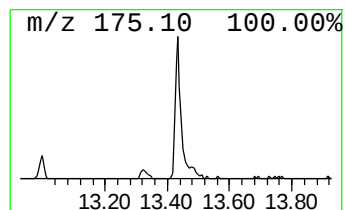
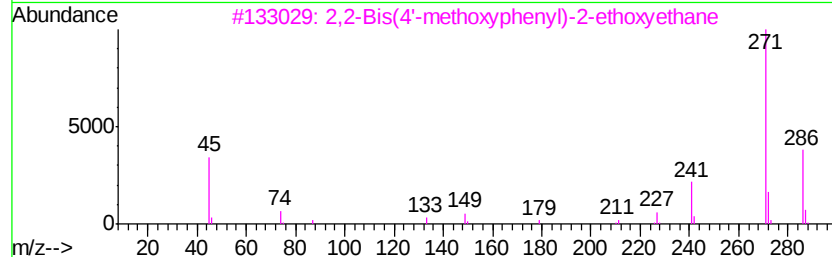
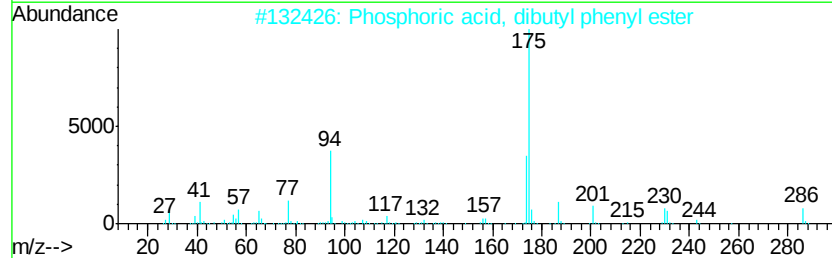
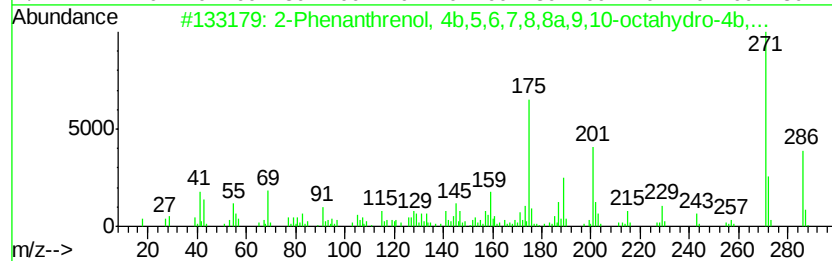
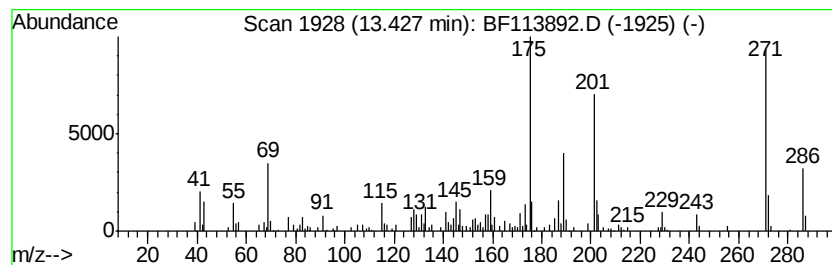
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.43	4.63 ng	221586	Chrysene-d12	14.05

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Phenanthrenol, 4b,5,6,7,8,8a,9...	286	C20H30O	000511-15-9	96
2		Phosphoric acid, dibutyl phenyl ...	286	C14H23O4P	002528-36-1	22
3		2,2-Bis(4'-methoxyphenyl)-2-etho...	286	C18H22O3	1000283-53-7	22
4		Benz(a)acridine, 9,10,12-trimethyl-	271	C20H17N	063040-02-8	18
5		6H-Indeno[1,2-d][1,2,4]-triazolo...	271	C16H9N5	1000286-52-6	14



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF042219\
Data File : BF113892.D
Acq On : 22 Apr 2019 21:50
Operator : JU/SJ
Sample : K2515-19
Misc :
ALS Vial : 10 Sample Multiplier: 1

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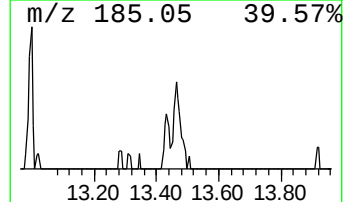
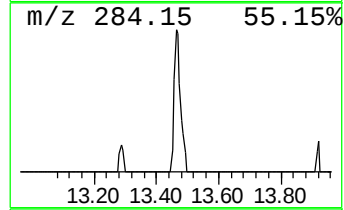
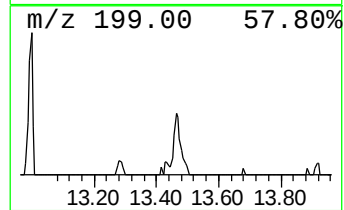
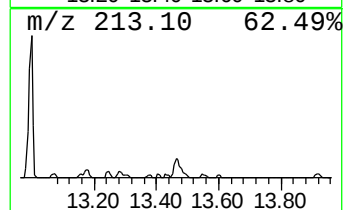
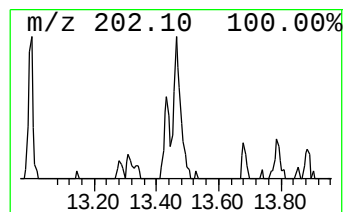
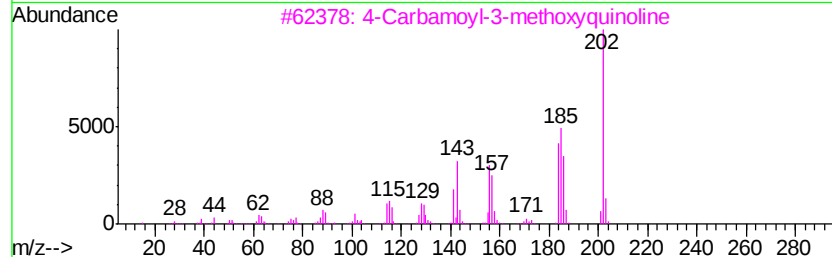
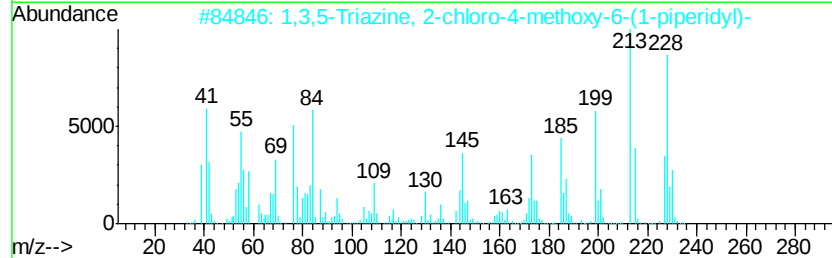
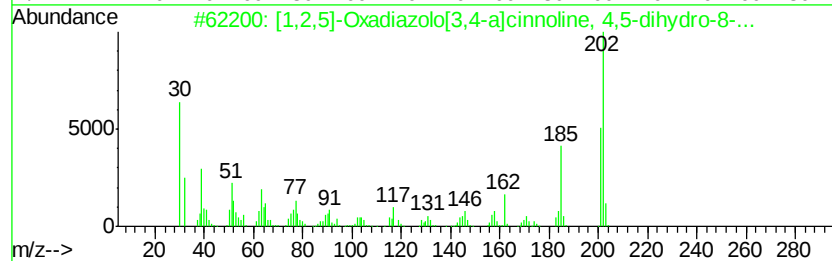
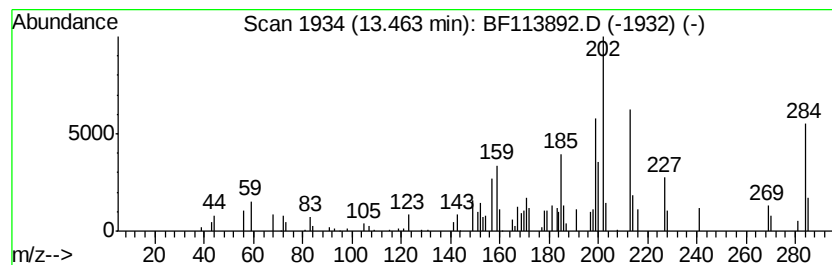
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 unknown13.46 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.46	2.46 ng	117732	Chrysene-d12	14.05

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	[1,2,5]-Oxadiazolo[3,4-a]cinnoli...	202	C10H10N4O	1000260-59-4	22
2		1,3,5-Triazine, 2-chloro-4-metho...	228	C9H13ClN4O	005097-48-3	22
3		4-Carbamoyl-3-methoxyquinoline	202	C11H10N2O2	020844-05-7	22
4		4-Methoxy-naphthalene-1-carboxyl...	202	C12H10O3	1000317-85-4	22
5		1-Naphthalenecarboxylic acid, 2-...	202	C12H10O3	000947-62-6	22



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF042219\
Data File : BF113892.D
Acq On : 22 Apr 2019 21:50
Operator : JU/SJ
Sample : K2515-19
Misc :
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Instrument :
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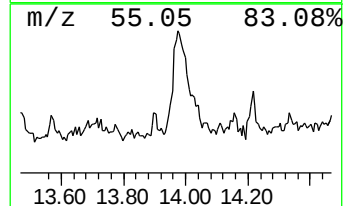
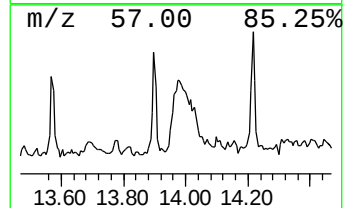
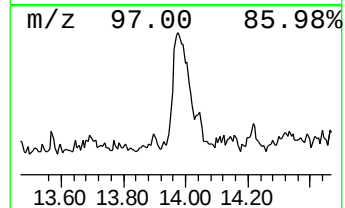
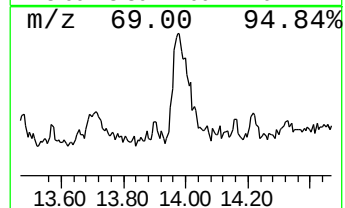
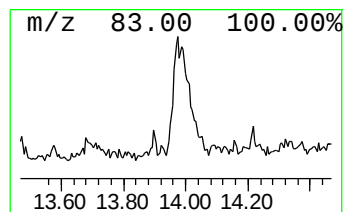
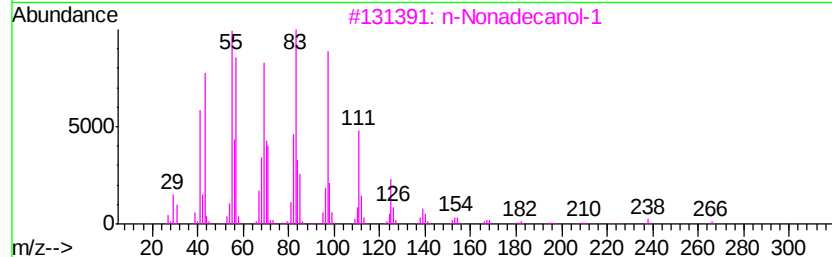
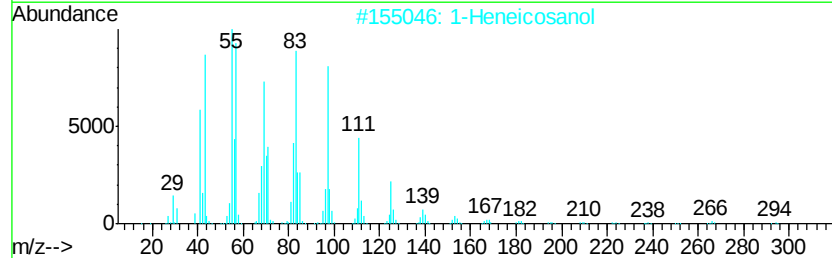
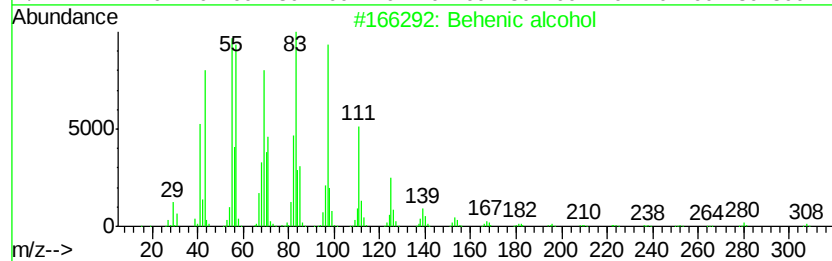
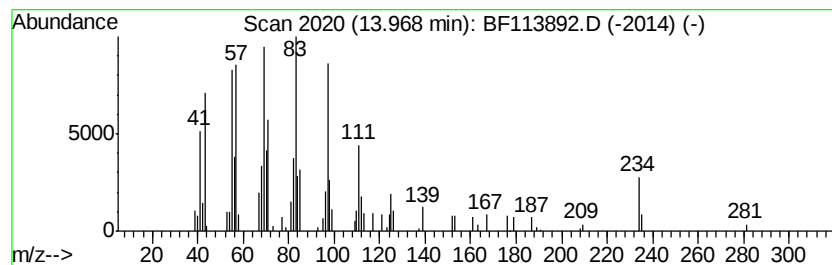
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Behenic alcohol Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.97	3.84 ng	184096	Chrysene-d12	14.05

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Behenic alcohol	326	C22H46O	000661-19-8	95
2		1-Heneicosanol	312	C21H44O	015594-90-8	95
3		n-Nonadecanol-1	284	C19H40O	001454-84-8	94
4		Pentafluoropropionic acid, penta...	374	C18H31F5O2	959092-08-1	93
5		Heptafluorobutyric acid, pentade...	424	C19H31F7O2	959261-23-5	93



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF042219\
Data File : BF113892.D
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Sample : K2515-19
Misc :
ALS Vial : 10 Sample Multiplier: 1

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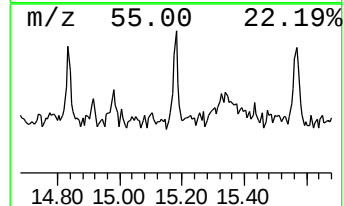
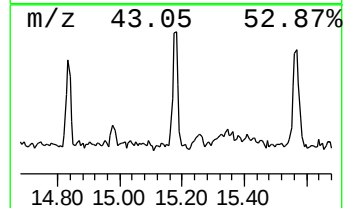
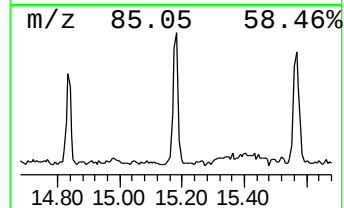
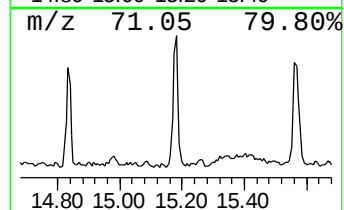
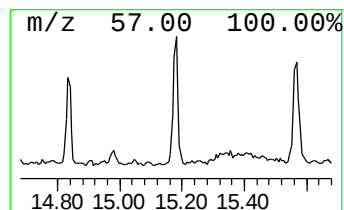
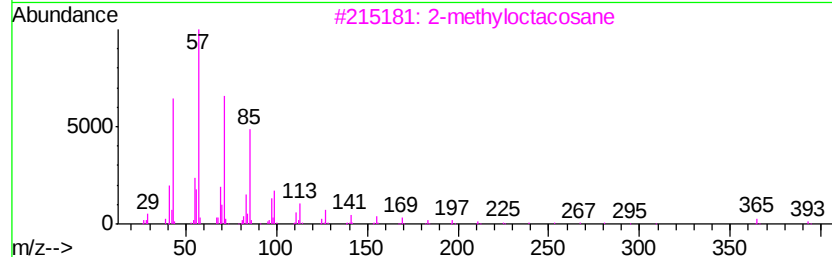
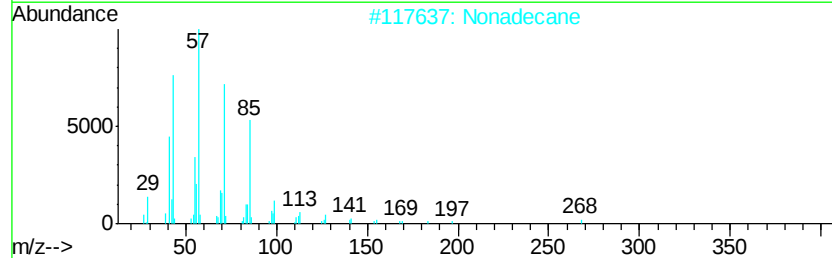
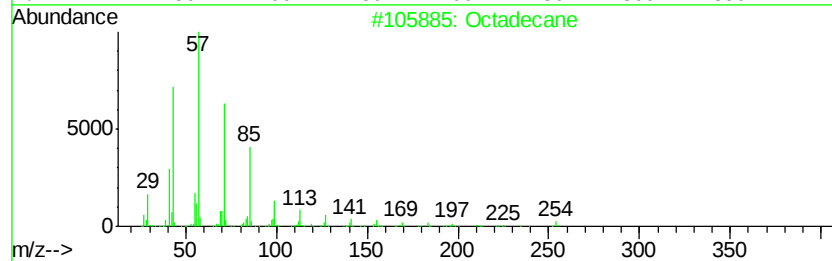
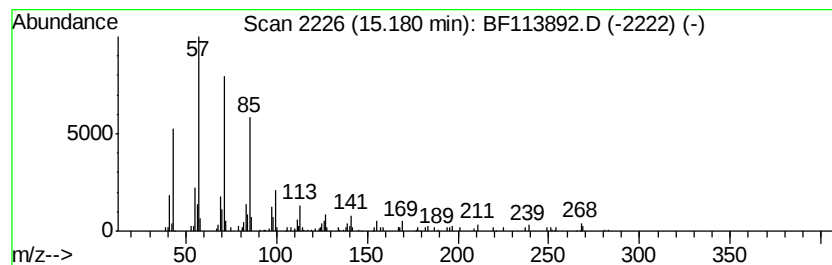
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Octadecane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.18	2.73 ng	133853	Perylene-d12	15.53

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecane	254	C18H38	000593-45-3	97
2		Nonadecane	268	C19H40	000629-92-5	93
3		2-methyloctacosane	408	C29H60	1000376-72-8	90
4		Triacontane	422	C30H62	000638-68-6	90
5		Octacosane	394	C28H58	000630-02-4	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF042219\
Data File : BF113892.D
Acq On : 22 Apr 2019 21:50
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Sample : K2515-19
Misc :
ALS Vial : 10 Sample Multiplier: 1

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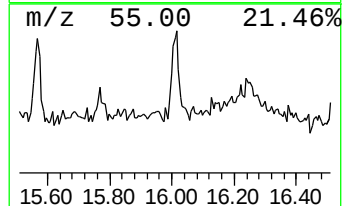
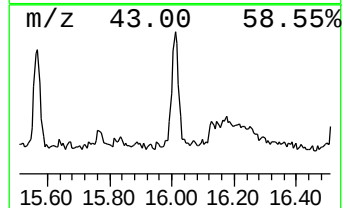
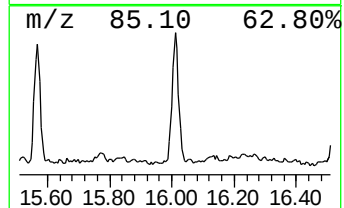
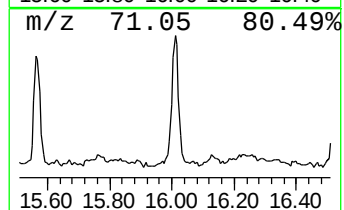
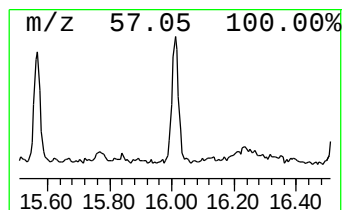
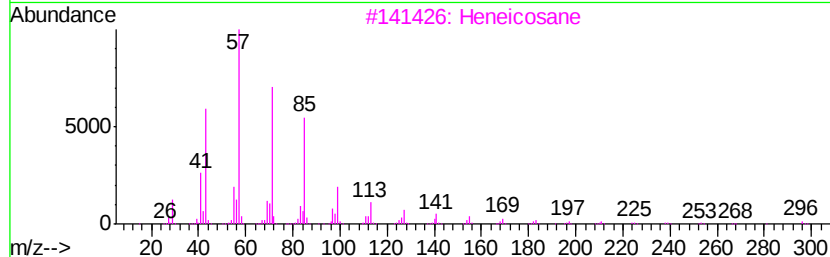
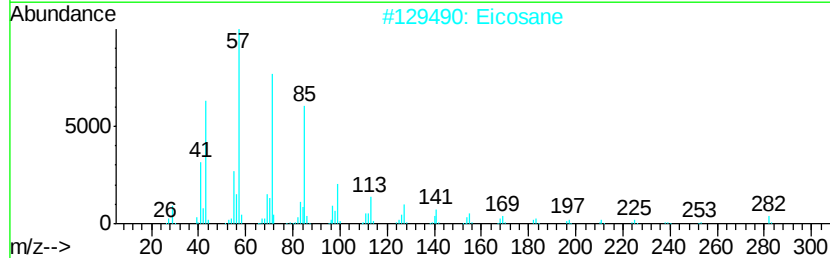
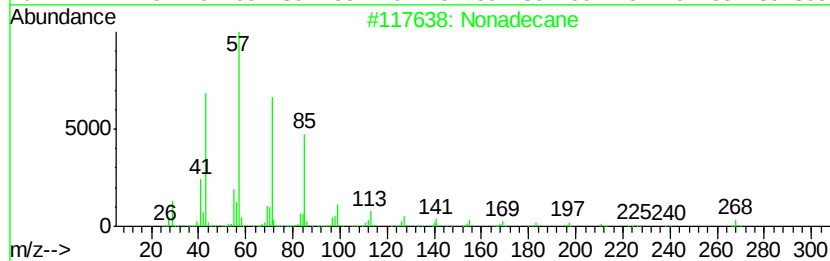
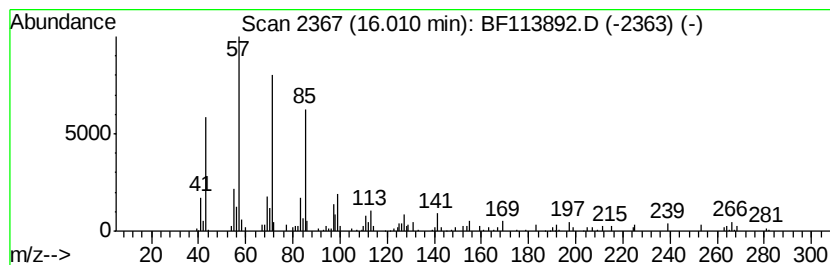
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Nonadecane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.01	3.04 ng	149134	Perylene-d12	15.53

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Nonadecane	268	C19H40	000629-92-5	95
2		Eicosane	282	C20H42	000112-95-8	93
3		Heneicosane	296	C21H44	000629-94-7	91
4		Octacosane	394	C28H58	000630-02-4	91
5		Nonadecane, 2,3-dimethyl-	296	C21H44	075163-99-4	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF042219\
Data File : BF113892.D
Acq On : 22 Apr 2019 21:50
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Sample : K2515-19
Misc :
ALS Vial : 10 Sample Multiplier: 1

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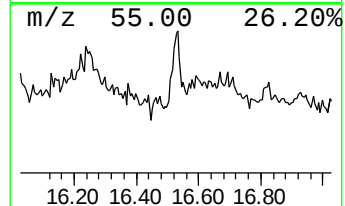
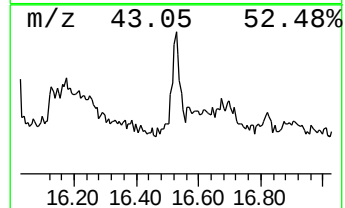
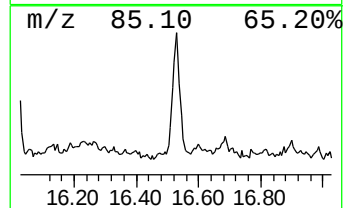
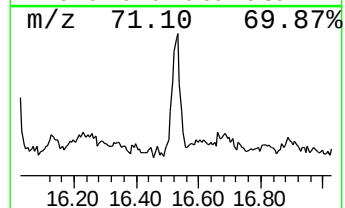
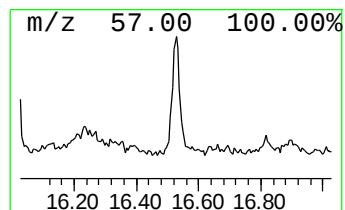
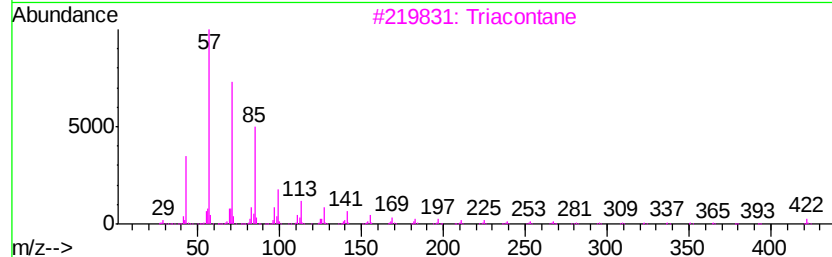
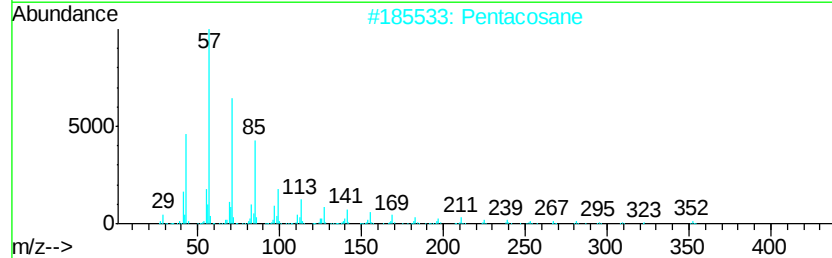
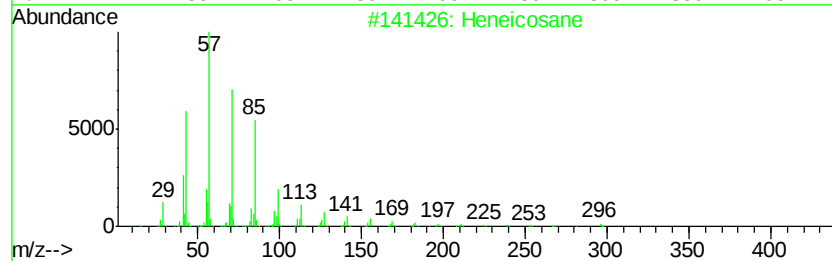
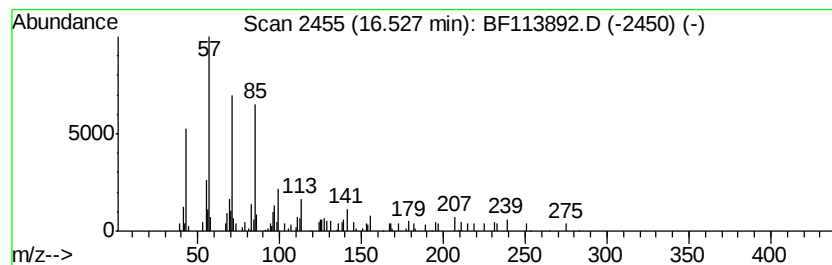
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Heneicosane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.53	2.11 ng	103364	Perylene-d12	15.53

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane		296	C21H44	000629-94-7	74
2	Pentacosane		352	C25H52	000629-99-2	74
3	Triacontane		422	C30H62	000638-68-6	72
4	Octacosane		394	C28H58	000630-02-4	72
5	2-methyloctacosane		408	C29H60	1000376-72-8	72



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF042219\
Data File : BF113892.D
Acq On : 22 Apr 2019 21:50
Operator : JU/SJ
Sample : K2515-19
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
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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
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unknown6.67	6.67	133.8	ng	4543980	1	6.89	679208	20.0
Longifolene	9.57	3.5	ng	152147	3	9.93	861851	20.0
n-Hexadecanoic acid	11.93	9.1	ng	357708	4	11.41	787192	20.0
Octadecanoic acid	12.72	4.7	ng	184150	4	11.41	787192	20.0
2-Phenanthrenol, ...	13.43	4.6	ng	221586	5	14.05	957965	20.0
unknown13.46	13.46	2.5	ng	117732	5	14.05	957965	20.0
Behenic alcohol	13.97	3.8	ng	184096	5	14.05	957965	20.0
Octadecane	15.18	2.7	ng	133853	6	15.53	981208	20.0
Nonadecane	16.01	3.0	ng	149134	6	15.53	981208	20.0
Heneicosane	16.53	2.1	ng	103364	6	15.53	981208	20.0