

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF110819\
 Data File : BF117740.D
 Acq On : 8 Nov 2019 19:23
 Operator : JU
 Sample : K5722-16
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 2-C

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-BF110619.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	2.234	19	25	33	rVB	1447446	1961347	24.26%	2.190%
2	5.151	515	521	527	rVB	1284503	1665002	20.59%	1.859%
3	5.545	582	588	591	rBV	6847254	6938494	85.82%	7.747%
4	6.551	752	759	762	rBV	6563871	7683458	95.04%	8.579%
5	6.698	779	784	787	rBV	7543496	7548939	93.37%	8.429%
6	6.928	819	823	826	rBV	2716412	2072340	25.63%	2.314%
7	7.086	846	850	853	rVB	6685104	5728264	70.85%	6.396%
8	7.486	913	918	921	rBV	4590270	4918987	60.84%	5.492%
9	8.216	1037	1042	1045	rBV	3246423	2801900	34.66%	3.128%
10	9.292	1220	1225	1229	rBV	8224456	8084572	100.00%	9.027%
11	9.375	1235	1239	1244	rBV4	135332	254761	3.15%	0.284%
12	9.686	1289	1292	1296	rBV	1297443	1320938	16.34%	1.475%
13	9.792	1307	1310	1313	rBV5	151604	150285	1.86%	0.168%
14	9.845	1315	1319	1322	rBV2	297724	482428	5.97%	0.539%
15	9.892	1324	1327	1330	rVB	455594	417037	5.16%	0.466%
16	9.974	1337	1341	1345	rVB	3498142	3390419	41.94%	3.785%
17	10.016	1346	1348	1349	rBV2	166913	139081	1.72%	0.155%
18	10.239	1383	1386	1388	rBV3	381659	444877	5.50%	0.497%
19	10.298	1393	1396	1398	rBV3	303729	414948	5.13%	0.463%
20	10.351	1403	1405	1408	rVV2	298281	311502	3.85%	0.348%
21	10.392	1409	1412	1416	rVV2	555001	645074	7.98%	0.720%
22	10.633	1450	1453	1456	rBV	1251911	1184946	14.66%	1.323%
23	10.769	1472	1476	1479	rBV	4970033	4753400	58.80%	5.307%
24	10.904	1495	1499	1502	rBV	3990887	4013290	49.64%	4.481%
25	11.374	1576	1579	1584	rVB	2997819	3067422	37.94%	3.425%
26	11.474	1593	1596	1598	rBV	2891375	2700529	33.40%	3.015%
27	13.068	1863	1867	1870	rVB	5147258	5053442	62.51%	5.642%
28	13.957	2016	2018	2026	rVB2	1514433	1761707	21.79%	1.967%
29	14.121	2043	2046	2049	rVB	2063788	1897463	23.47%	2.119%
30	14.986	2189	2193	2201	rVB	1288550	1461422	18.08%	1.632%
31	15.174	2222	2225	2229	rBV	777085	1163419	14.39%	1.299%
32	15.256	2237	2239	2242	rVB	221071	210133	2.60%	0.235%
33	15.492	2275	2279	2285	rVB	488420	609569	7.54%	0.681%
34	15.556	2287	2290	2294	rVV	384391	446572	5.52%	0.499%

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

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

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Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	15.621	2297	2301	2305	rVV	1673155	2225424	27.53%	2.485%
36	15.656	2305	2307	2314	rVB3	342351	468256	5.79%	0.523%
37	16.121	2383	2386	2391	rVB	138563	179762	2.22%	0.201%
38	17.139	2554	2559	2567	rVB2	246637	504630	6.24%	0.563%
39	17.603	2633	2638	2654	rVB	211612	488208	6.04%	0.545%

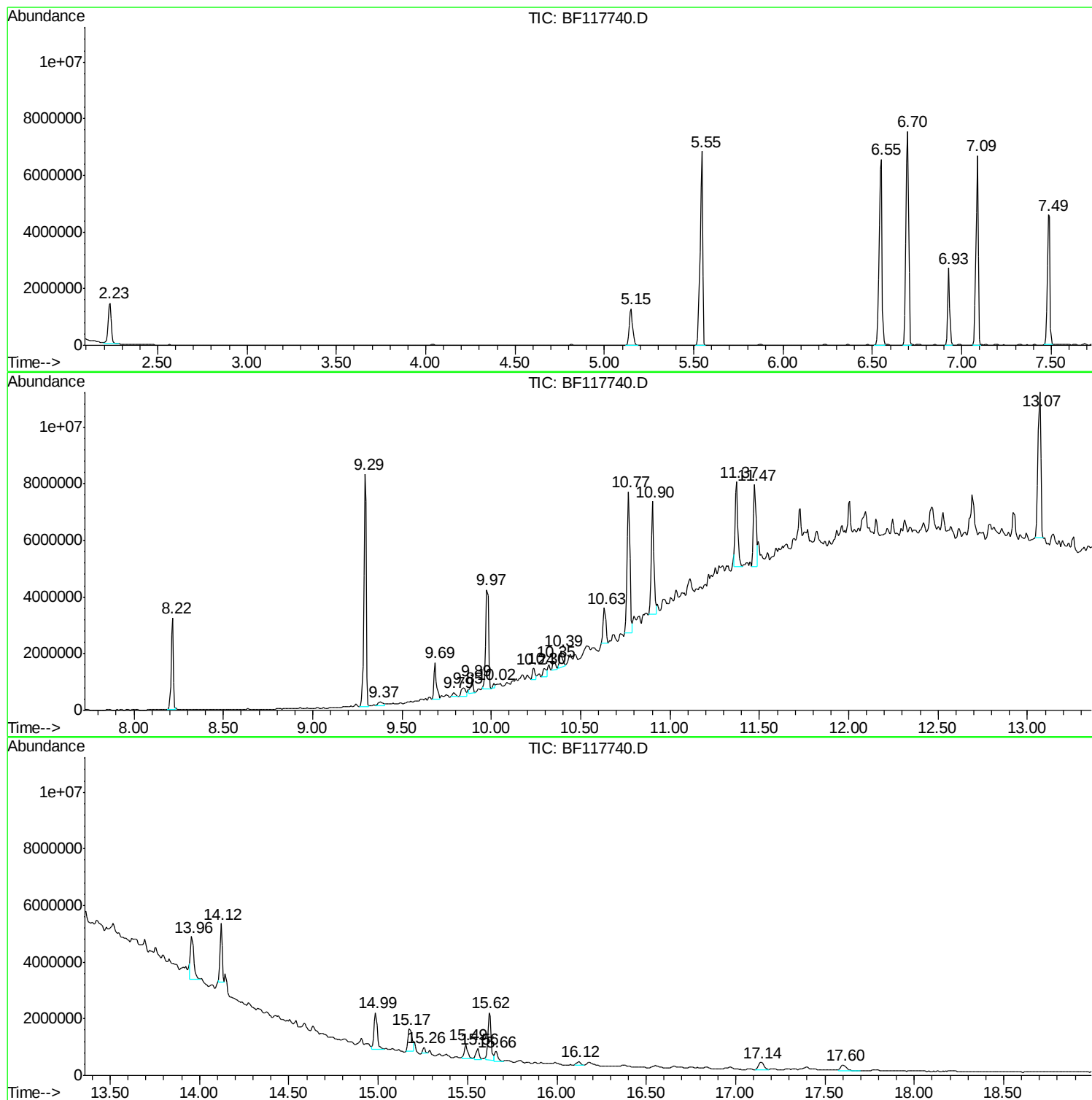
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF110619.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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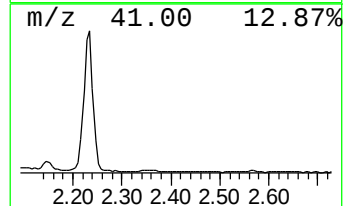
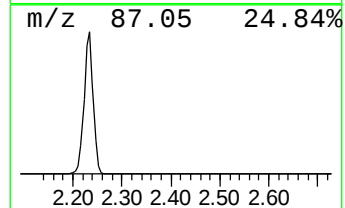
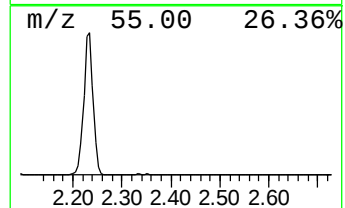
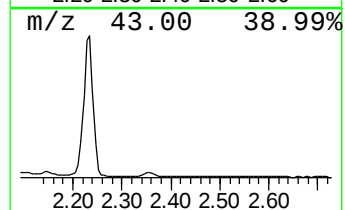
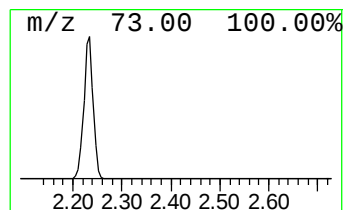
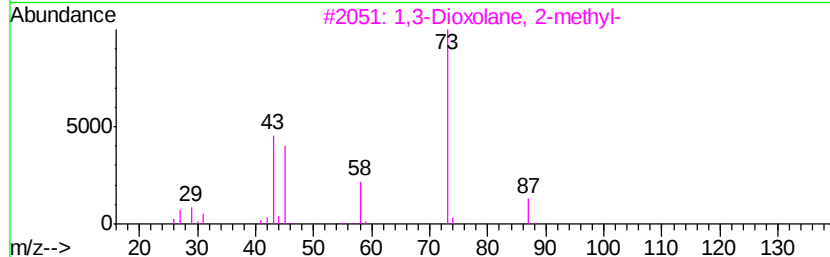
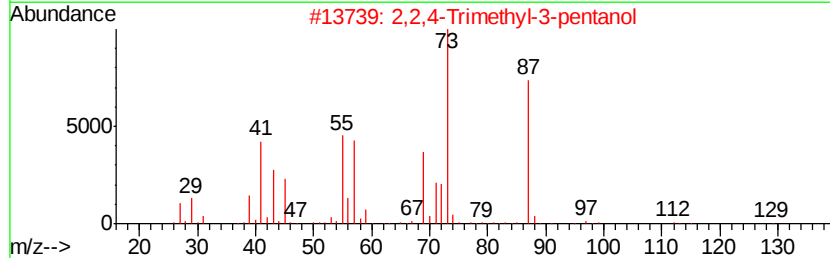
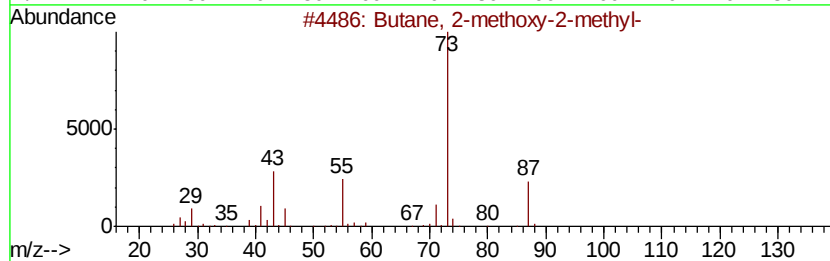
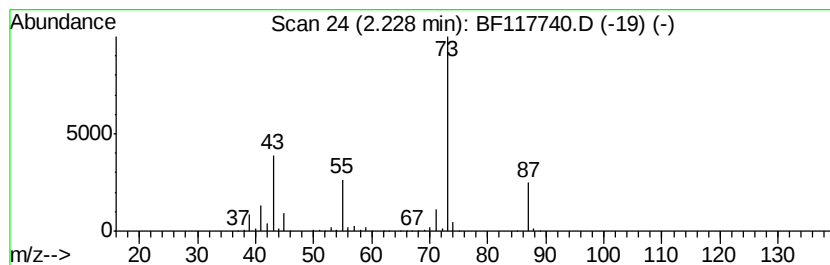
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF110619.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.23	18.93 ng	1961350	1,4-Dichlorobenzene-d4	6.93

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		1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	37
4		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	12
5		Silane, tetramethyl-	88	C4H12Si	000075-76-3	10



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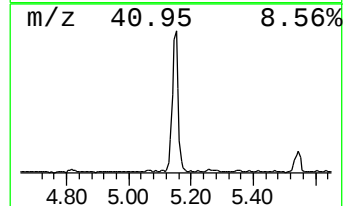
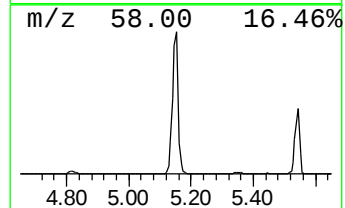
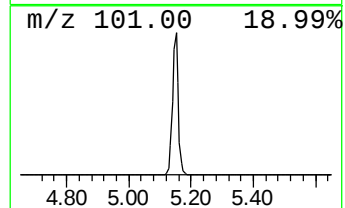
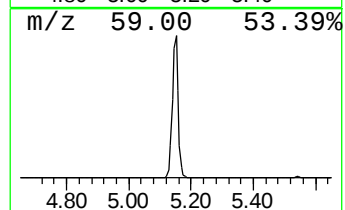
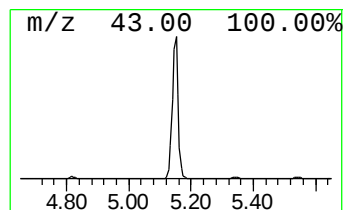
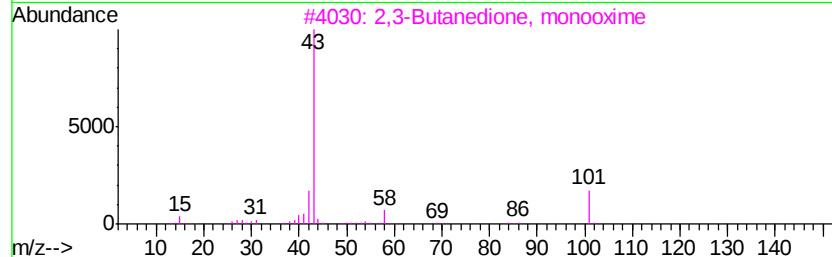
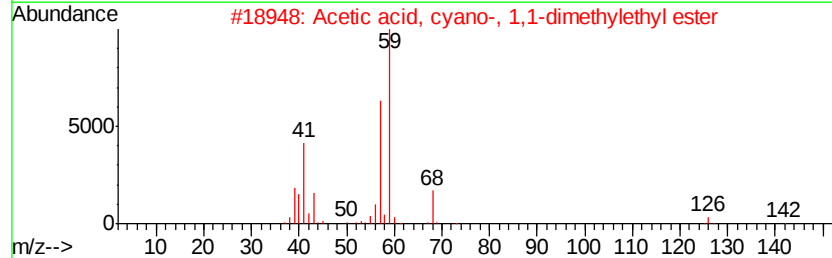
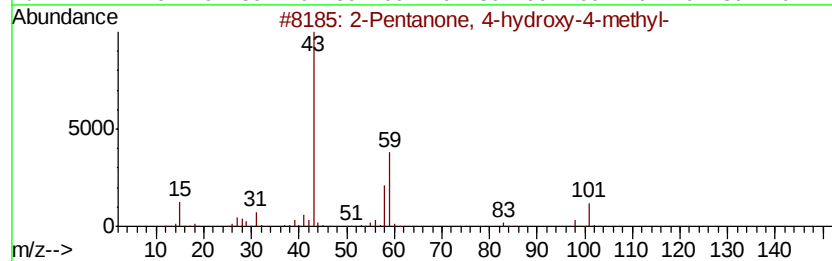
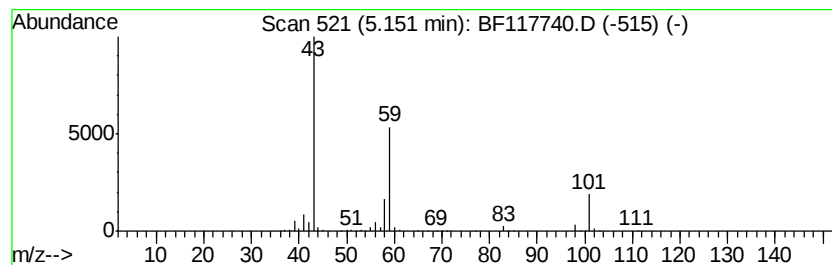
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Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.15	16.07 ng	1665000	1,4-Dichlorobenzene-d4	6.93

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, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	25
3		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	16
4		1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10
5		Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9



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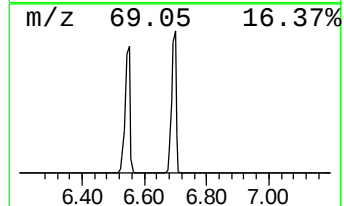
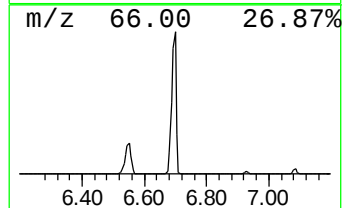
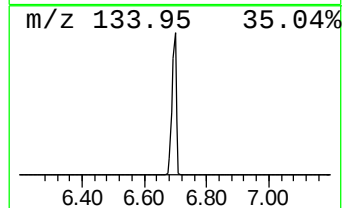
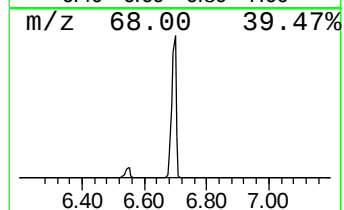
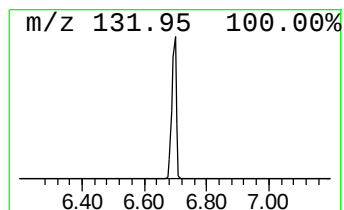
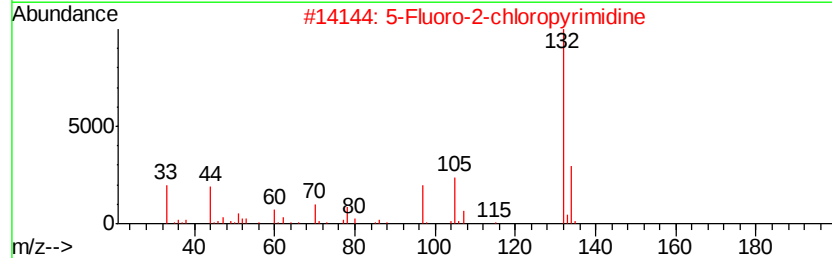
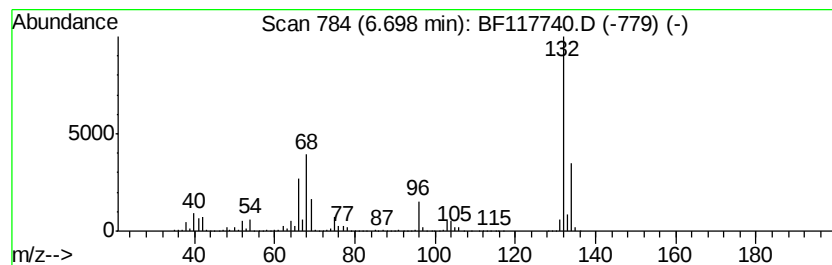
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 3 unknown6.70 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.70	72.85 ng	7548940	1,4-Dichlorobenzene-d4	6.93

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16
3		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	12
4		5-Aminoindole	132	C8H8N2	005192-03-0	12
5		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9



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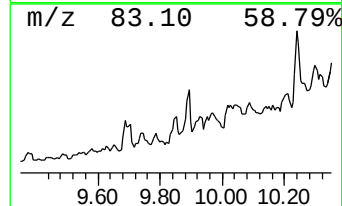
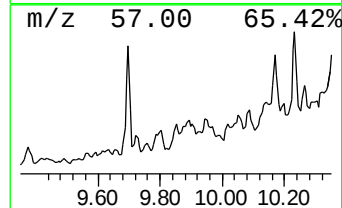
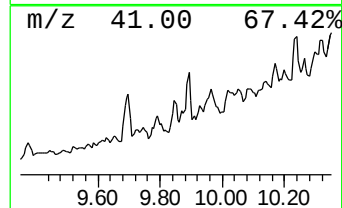
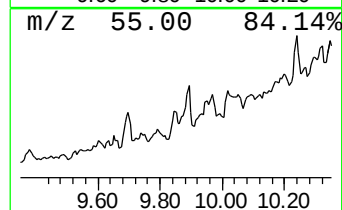
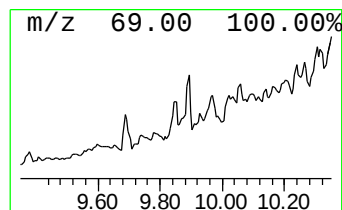
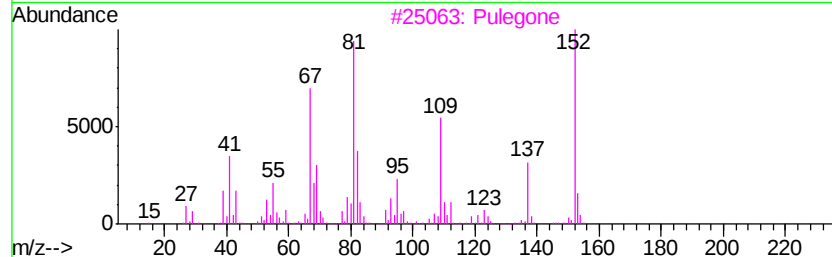
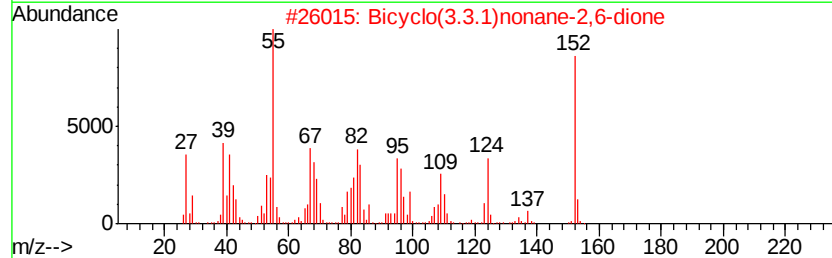
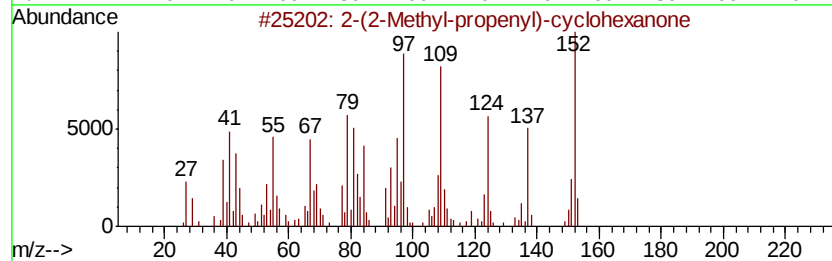
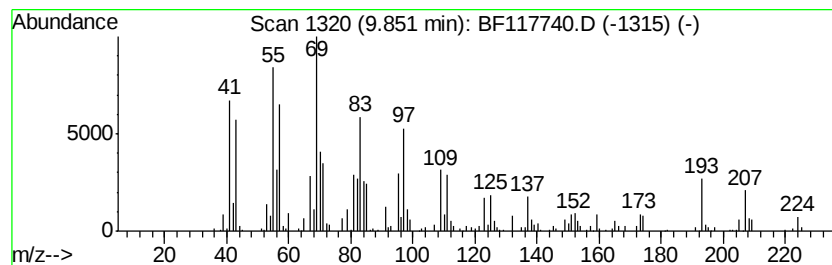
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 5 unknown9.85 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.85	2.85 ng	482428	Acenaphthene-d10	9.97

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-(2-Methyl-propenyl)-cyclohexanone	152	C10H16O	065737-44-2	45
2		Bicyclo(3.3.1)nonane-2,6-dione	152	C9H12O2	016473-11-3	25
3		Pulegone	152	C10H16O	000089-82-7	25
4		Trimethyl thioborate	152	C3H9BS3	000997-49-9	25
5		Bicyclo[3.1.1]heptan-2-one, 3,6,...	152	C10H16O	016022-08-5	25



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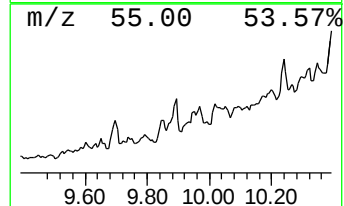
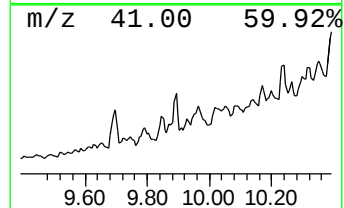
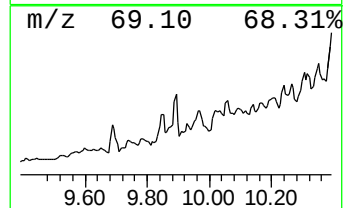
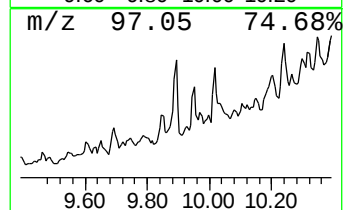
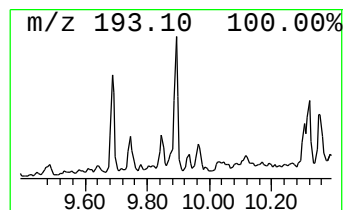
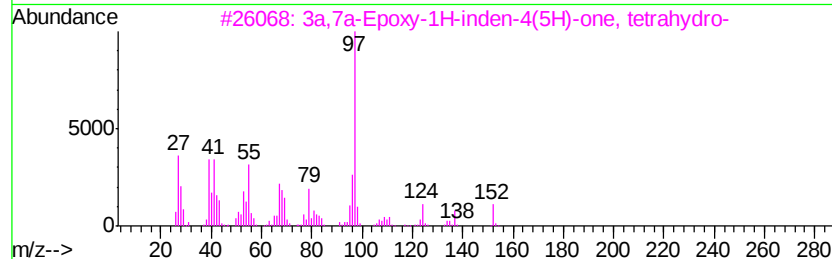
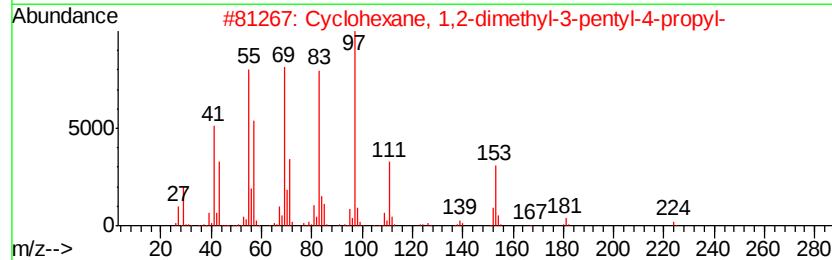
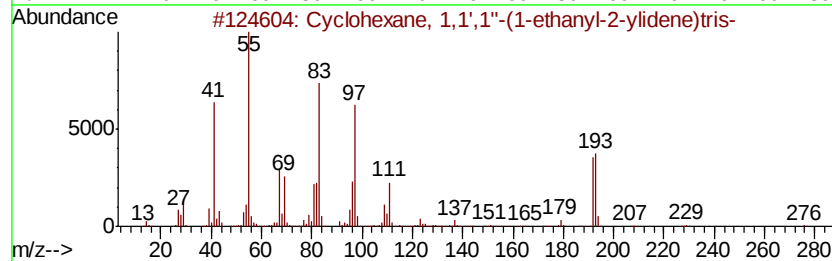
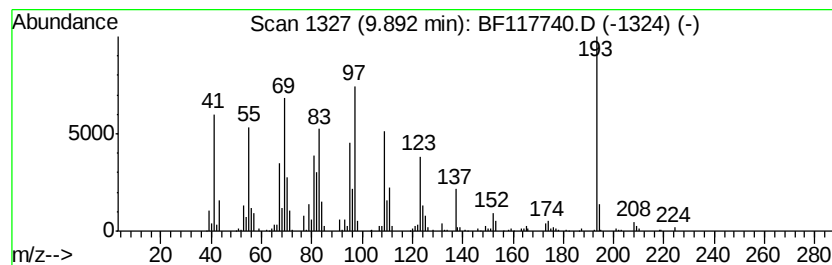
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Peak Number 6 unknown9.89 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.89	2.46 ng	417037	Acenaphthene-d10	9.97

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexane, 1,1',1''-(1-ethanyl...	276	C20H36	055682-86-5	43
2		Cyclohexane, 1,2-dimethyl-3-pent...	224	C16H32	062376-17-4	38
3		3a,7a-Epoxy-1H-inden-4(5H)-one, ...	152	C9H12O2	039746-31-1	18
4		Trans-1-methyl-2-nonyl-cyclohexane	224	C16H32	1000371-47-8	18
5		Aziridine, 1-(2-methyl-1-propenyl)-	97	C6H11N	080839-93-6	15



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF110819\
Data File : BF117740.D
Acq On : 8 Nov 2019 19:23
Operator : JU
Sample : K5722-16
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
2-C

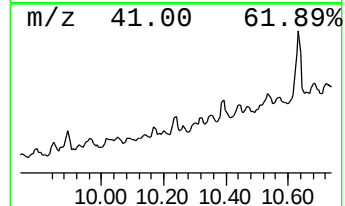
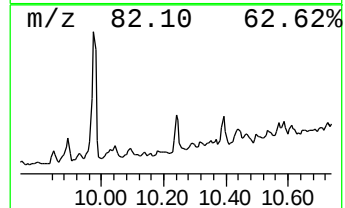
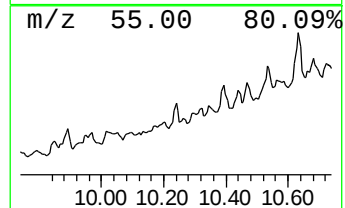
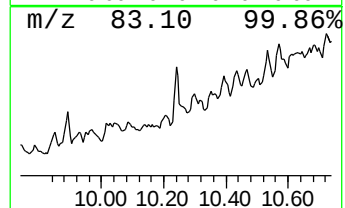
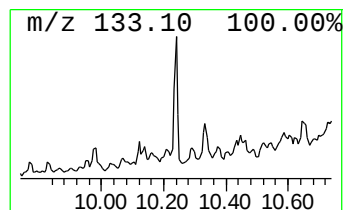
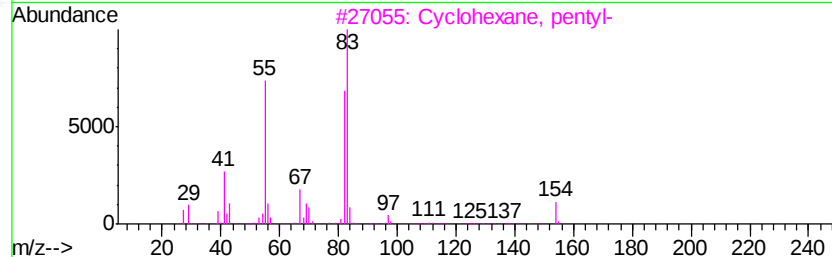
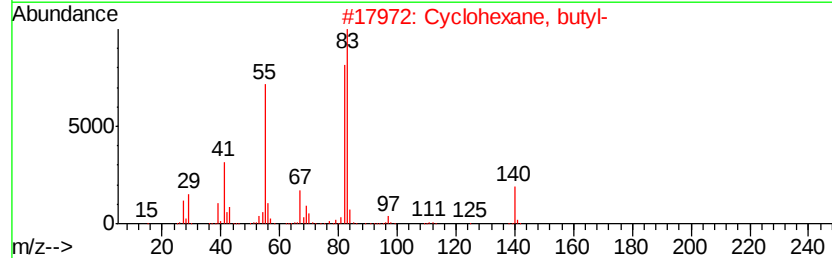
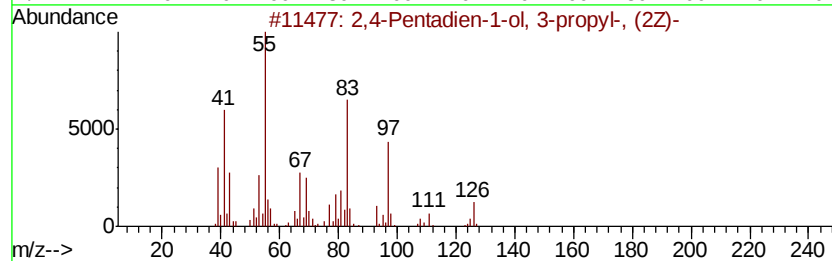
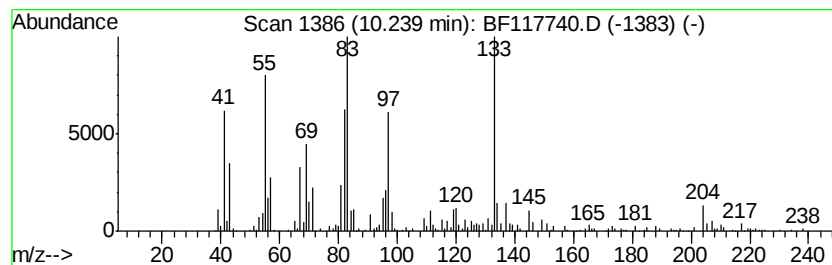
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 unknown10.24 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.24	2.62 ng	444877	Acenaphthene-d10	9.97

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2,4-Pentadien-1-ol, 3-propyl-, (...)	126	C8H14O	1000142-17-9	38
2		Cyclohexane, butyl-	140	C10H20	001678-93-9	27
3		Cyclohexane, pentyl-	154	C11H22	004292-92-6	27
4		Cyclohexane, propyl-	126	C9H18	001678-92-8	27
5		Cyclohexane, ethyl-	112	C8H16	001678-91-7	27



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF110819\
Data File : BF117740.D
Acq On : 8 Nov 2019 19:23
Operator : JU
Sample : K5722-16
Misc :
ALS Vial : 13 Sample Multiplier: 1

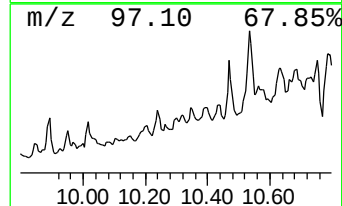
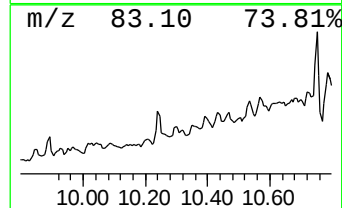
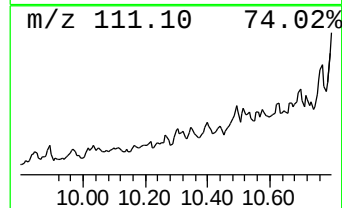
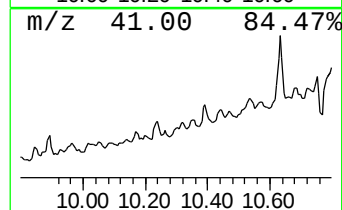
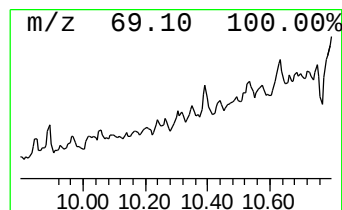
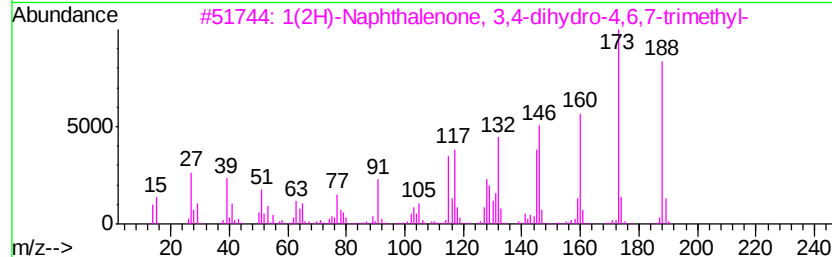
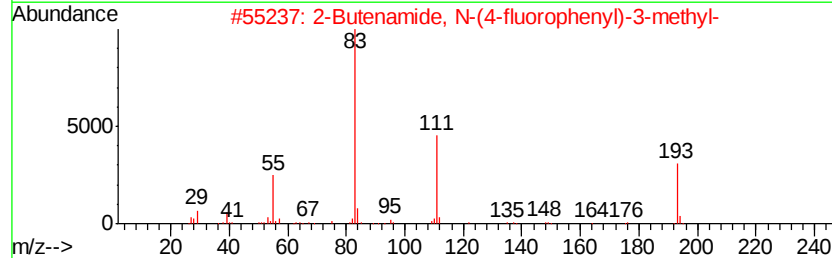
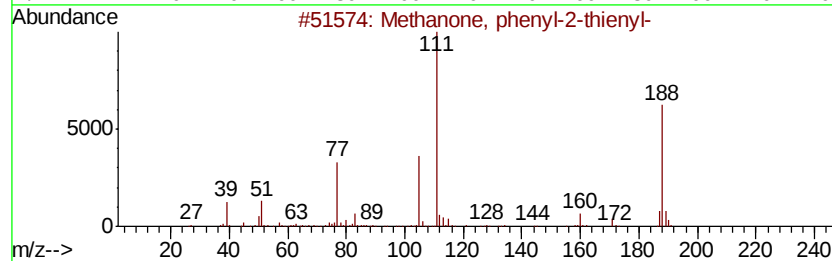
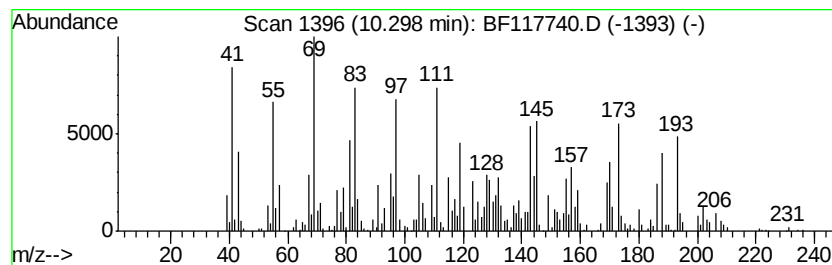
Instrument :
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF110619.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 unknown10.30 Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.30	2.45 ng	414948	Acenaphthene-d10	9.97
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Methanone, phenyl-2-thienyl-	188 C11H8OS	000135-00-2 14
2		2-Butenamide, N-(4-fluorophenyl)...	193 C11H12FNO	1000307-26-7 14
3		1(2H)-Naphthalenone, 3,4-dihydro...	188 C13H16O	030316-32-6 11
4		Benzenamine, 2-fluoro-	111 C6H6FN	000348-54-9 11
5		Cyclohexanone, 3-(4-hydroxybutyl...	184 C11H20O2	091212-98-5 11



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF110819\
Data File : BF117740.D
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Sample : K5722-16
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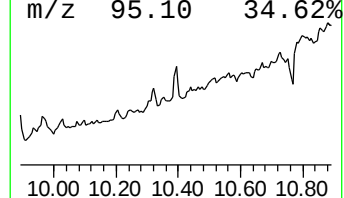
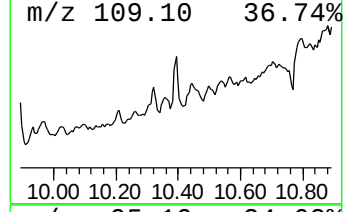
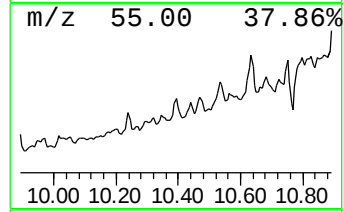
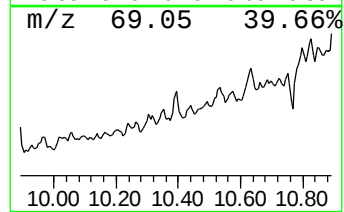
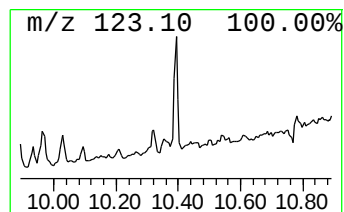
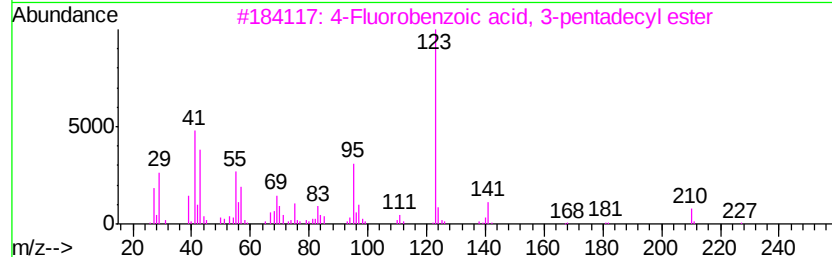
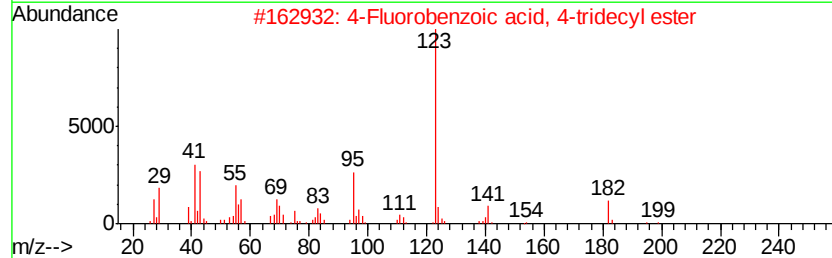
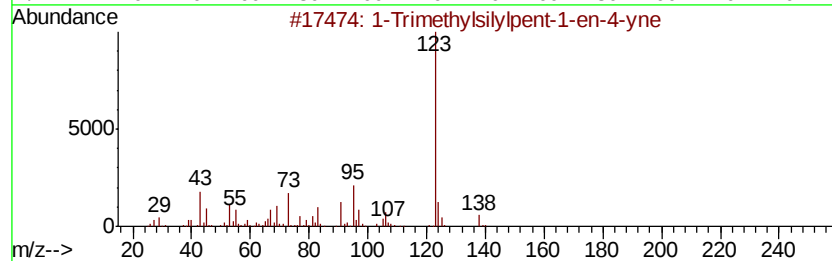
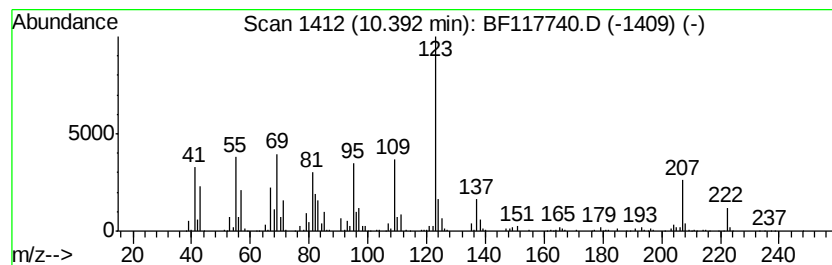
Instrument :
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF110619.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 1-Trimethylsilylpent-1-en-4... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.39	3.81 ng	645074	Acenaphthene-d10	9.97	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Trimethylsilylpent-1-en-4-yne	138	C8H14Si	079516-25-9	50
2	4-Fluorobenzoic acid, 4-tridecyl...	322	C20H31F02	1000280-68-0	50
3	4-Fluorobenzoic acid, 3-pentadec...	350	C22H35F02	1000280-68-4	50
4	2-Fluorobenzoic acid, oct-3-en-2...	250	C15H19F02	1000299-16-5	50
5	3-Fluorobenzoic acid, 2-ethylcyc...	250	C15H19F02	1000279-04-9	47



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF110819\
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Sample : K5722-16
Misc :
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Instrument :
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ClientSampleId :
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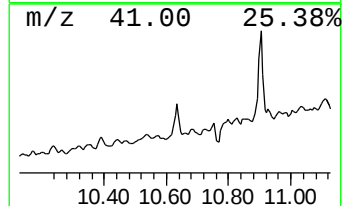
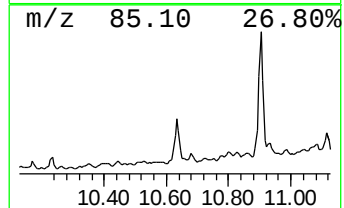
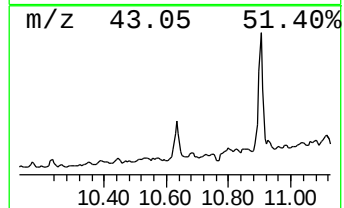
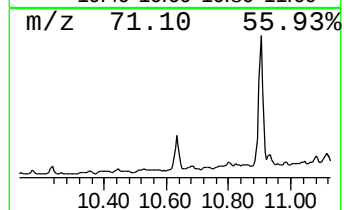
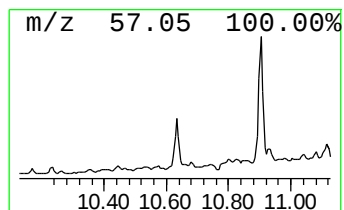
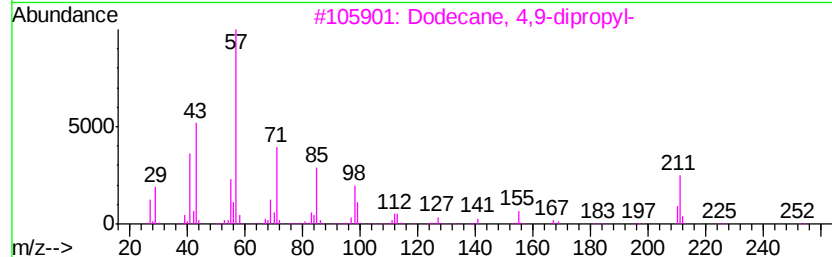
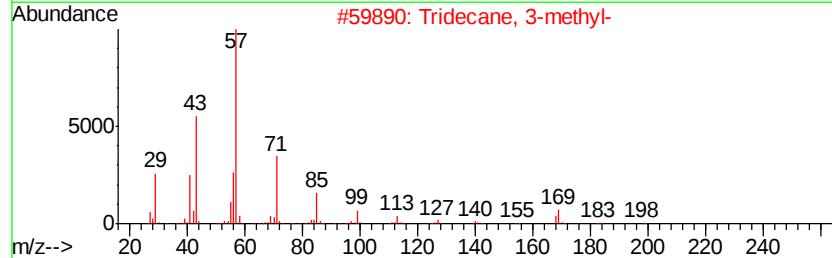
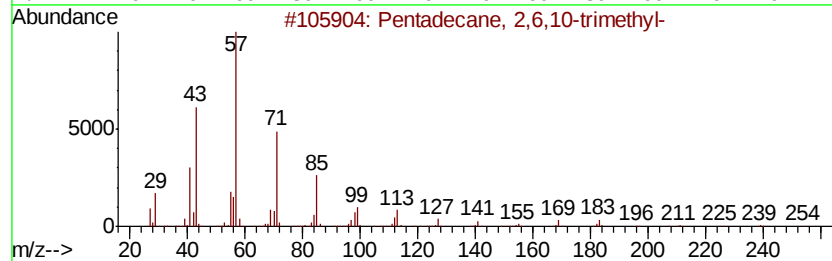
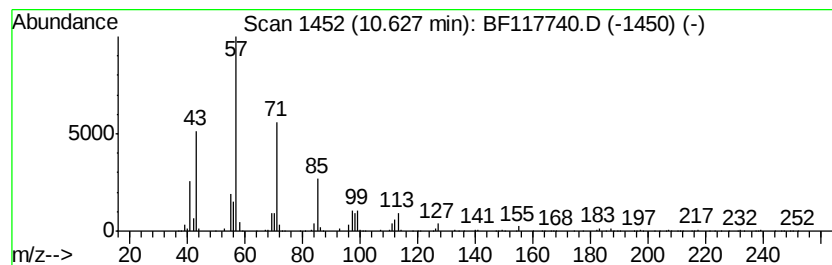
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Pentadecane, 2,6,10-trimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.63	6.99 ng	1184950	Acenaphthene-d10	9.97

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentadecane, 2,6,10-trimethyl-	254	C18H38	003892-00-0	91
2		Tridecane, 3-methyl-	198	C14H30	006418-41-3	90
3		Dodecane, 4,9-dipropyl-	254	C18H38	003054-63-5	87
4		Heptacosane	380	C27H56	000593-49-7	86
5		Tridecane, 6-methyl-	198	C14H30	013287-21-3	72



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF110819\
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Sample : K5722-16
Misc :
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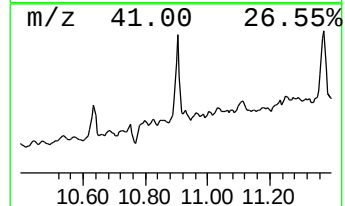
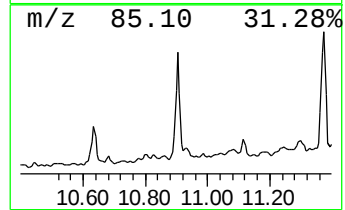
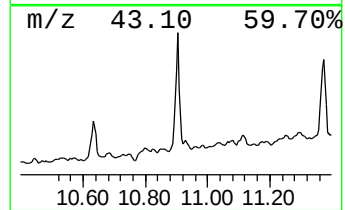
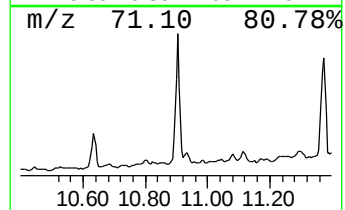
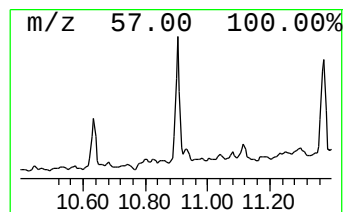
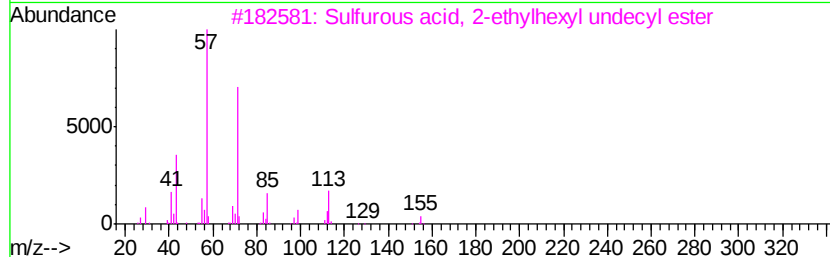
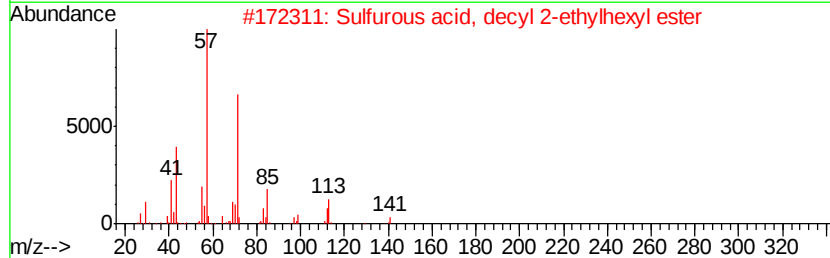
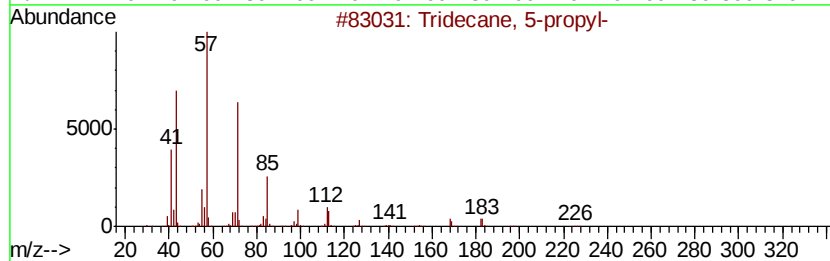
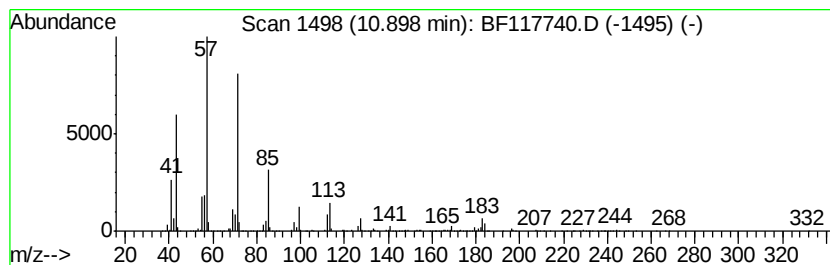
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Tridecane, 5-propyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.90	29.72 ng	4013290	Phenanthrene-d10	11.47	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tridecane, 5-propyl-	226	C16H34	055045-11-9	83
2	Sulfurous acid, decyl 2-ethylhex...	334	C18H38O3S	1000309-19-3	72
3	Sulfurous acid, 2-ethylhexyl und...	348	C19H40O3S	1000309-19-4	72
4	Dodecane, 2,7,10-trimethyl-	212	C15H32	074645-98-0	72
5	Undecane, 6-ethyl-	184	C13H28	017312-60-6	72



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF110819\
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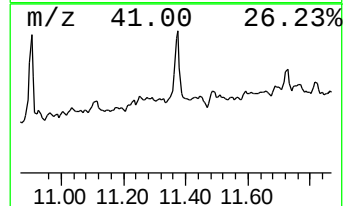
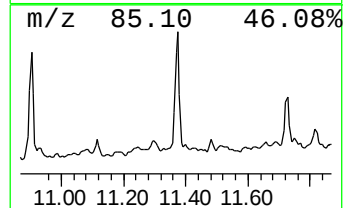
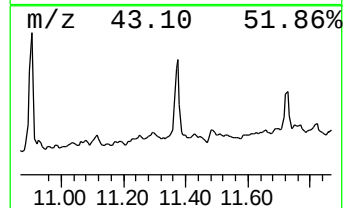
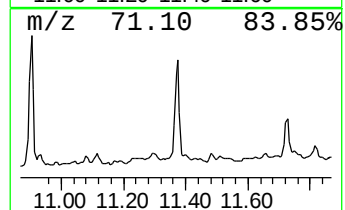
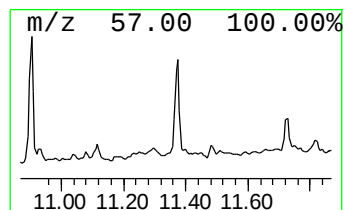
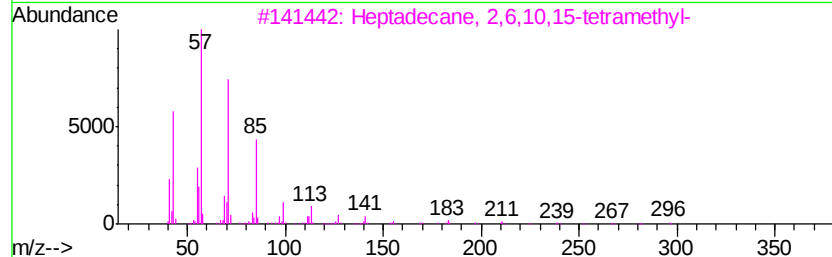
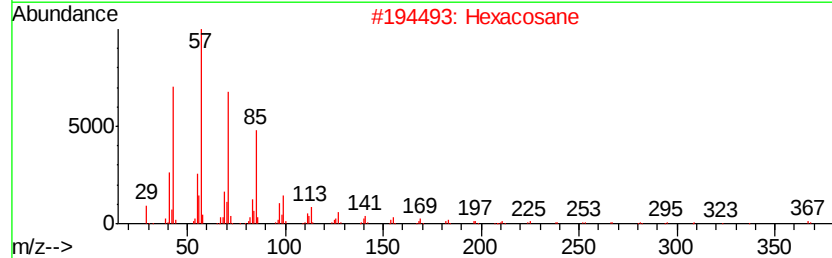
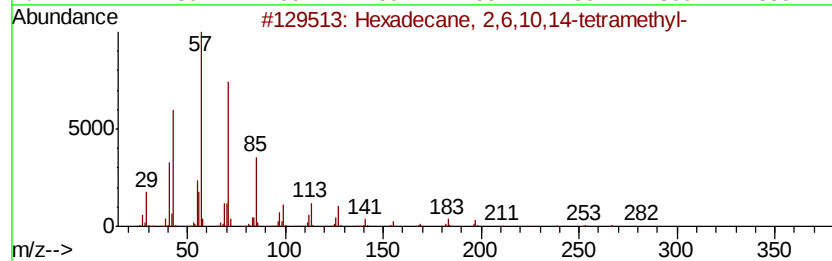
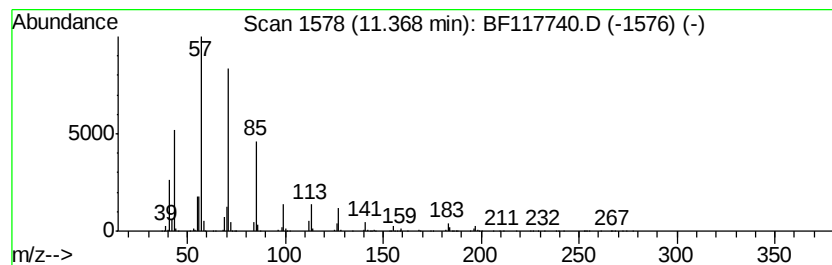
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

 Peak Number 12 Hexadecane, 2,6,10,14-tetra... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.37	22.72 ng	3067420	Phenanthrene-d10	11.47

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadecane, 2,6,10,14-tetramethyl-	282	C20H42	000638-36-8	91
2		Hexacosane	366	C26H54	000630-01-3	90
3		Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	86
4		Pentadecane, 4-methyl-	226	C16H34	002801-87-8	86
5		Dodecane, 4-methyl-	184	C13H28	006117-97-1	74



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF110819\
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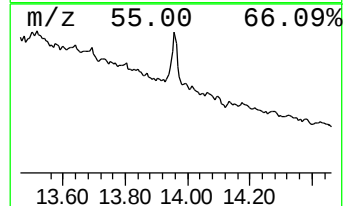
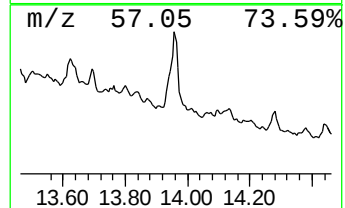
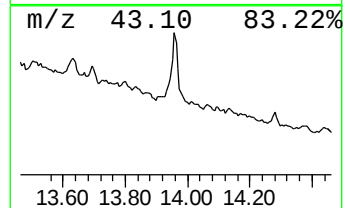
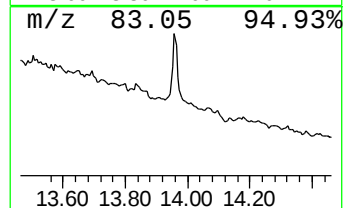
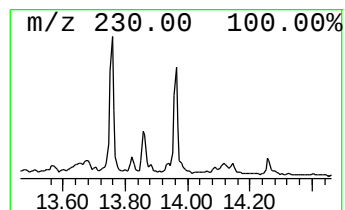
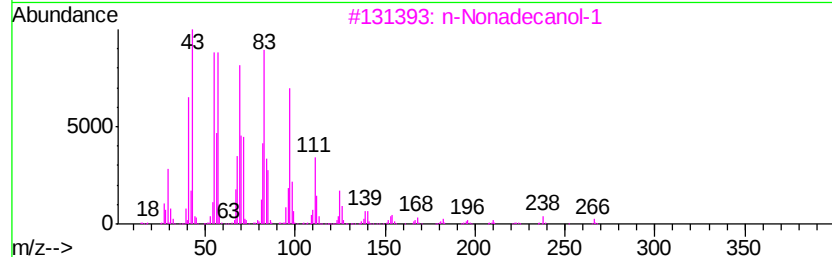
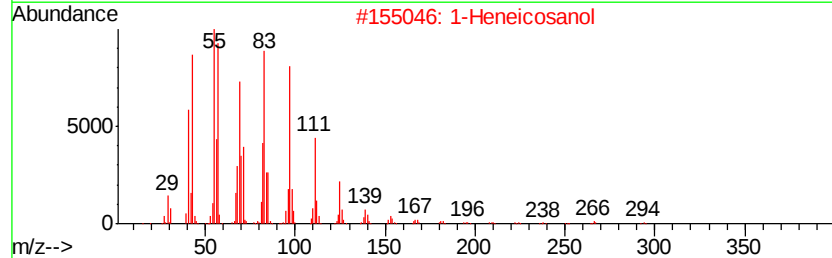
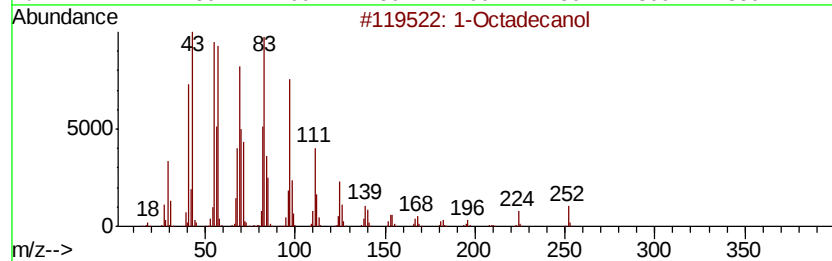
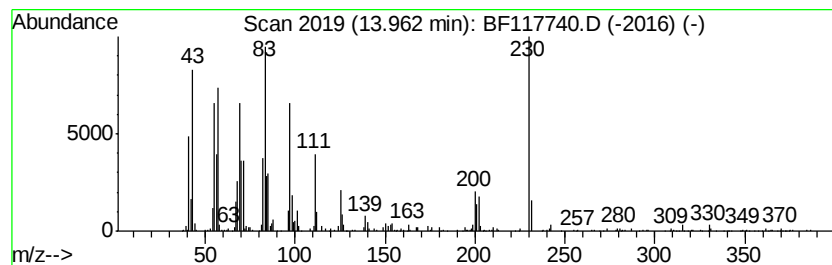
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 1-Octadecanol Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.96	18.57 ng	1761710	Chrysene-d12	14.12

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Octadecanol	270	C18H38O	000112-92-5	93
2		1-Heneicosanol	312	C21H44O	015594-90-8	93
3		n-Nonadecanol-1	284	C19H40O	001454-84-8	93
4		Trifluoroacetic acid, pentadecyl...	324	C17H31F3O2	959010-23-2	89
5		n-Heptadecanol-1	256	C17H36O	001454-85-9	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF110819\
Data File : BF117740.D
Acq On : 8 Nov 2019 19:23
Operator : JU
Sample : K5722-16
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
2-C

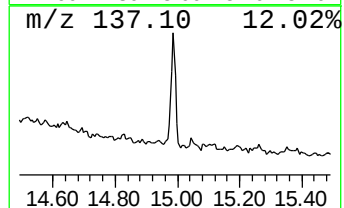
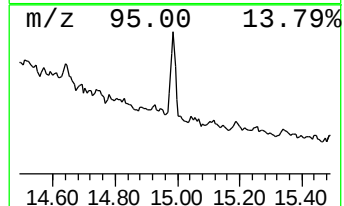
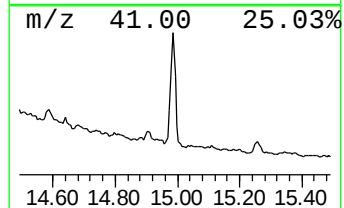
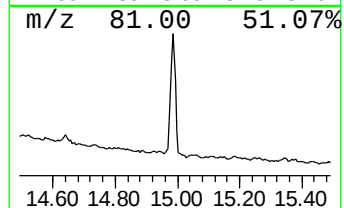
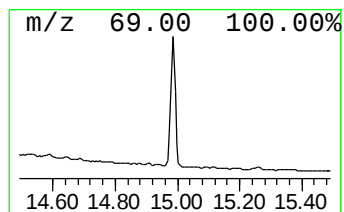
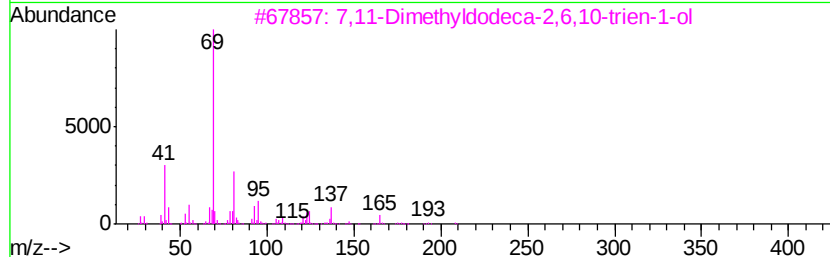
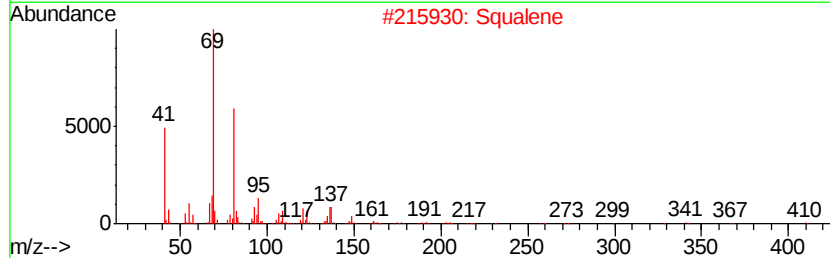
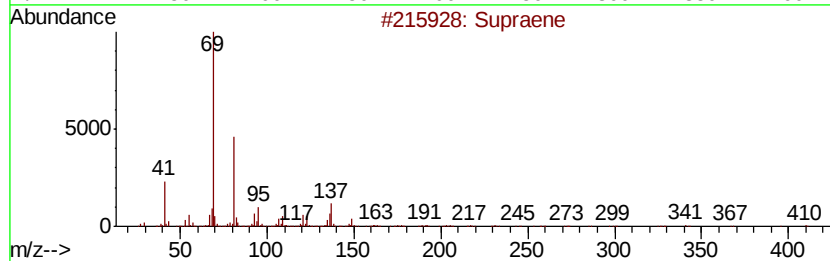
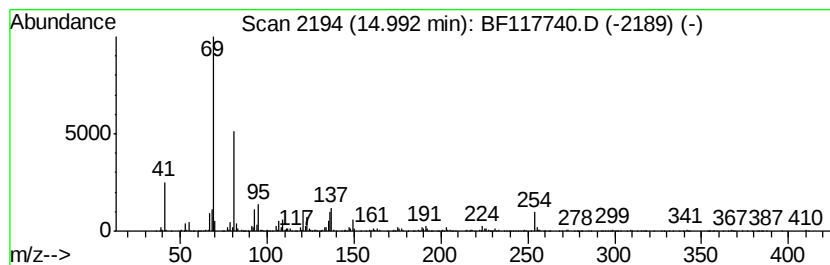
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Supraene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.99	13.13 ng	1461420	Perylene-d12	15.62

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Supraene	410	C30H50	007683-64-9	91
2		Squalene	410	C30H50	000111-02-4	91
3		7,11-Dimethyldodeca-2,6,10-trien...	208	C14H24O	125529-10-4	64
4		.beta.-Citronellol, chlorodifluo...	268	C12H19ClF2O2	1000376-20-8	53
5		1,5-Heptadiene, 3,3,6-trimethyl-	138	C10H18	035387-63-4	52



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF110819\
Data File : BF117740.D
Acq On : 8 Nov 2019 19:23
Operator : JU
Sample : K5722-16
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
2-C

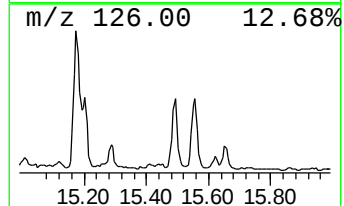
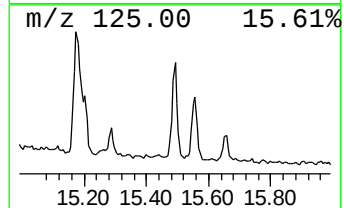
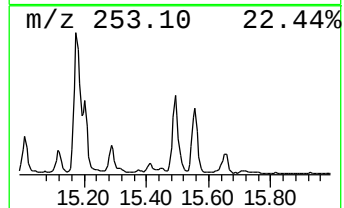
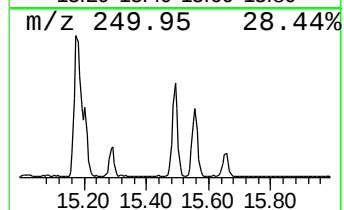
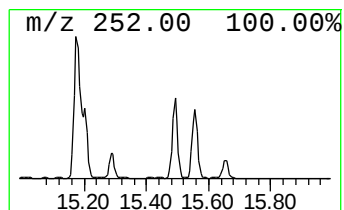
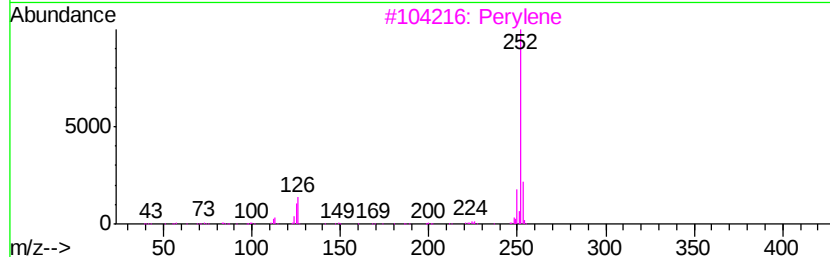
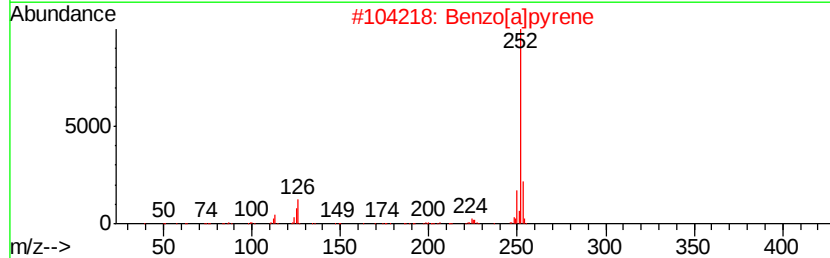
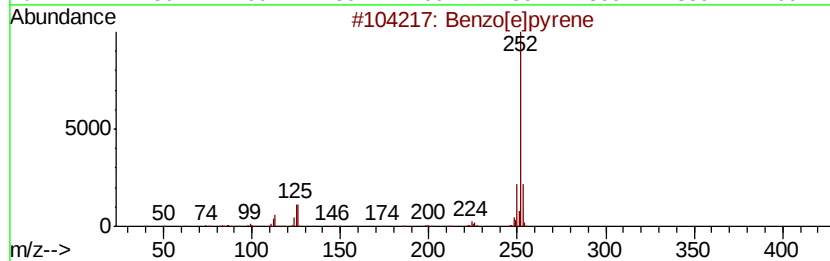
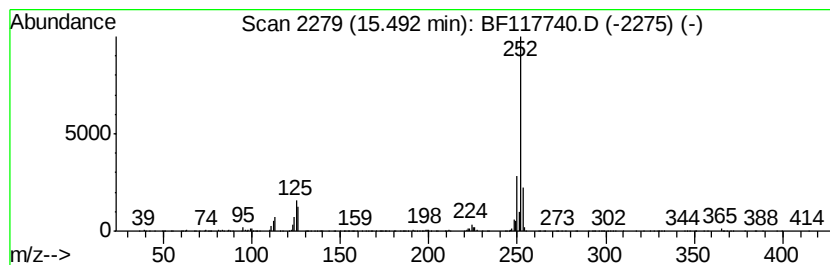
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Benzo[e]pyrene Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.49	5.48 ng	609569	Perylene-d12	15.62

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo[e]pyrene	252	C20H12	000192-97-2	99
2			Benzo[a]pyrene	252	C20H12	000050-32-8	97
3			Perylene	252	C20H12	000198-55-0	94
4			Benzo[e]acephenanthrylene	252	C20H12	000205-99-2	93
5			Benzo[k]fluoranthene	252	C20H12	000207-08-9	93



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF110819\
 Data File : BF117740.D
 Acq On : 8 Nov 2019 19:23
 Operator : JU
 Sample : K5722-16
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 2-C

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF110619.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...	2.23	18.9	ng	1961350	1	6.93	2072340	20.0
2-Pentanone, 4-hy...	5.15	16.1	ng	1665000	1	6.93	2072340	20.0
unknown6.70	6.70	72.8	ng	7548940	1	6.93	2072340	20.0
unknown9.85	9.85	2.9	ng	482428	3	9.97	3390420	20.0
unknown9.89	9.89	2.5	ng	417037	3	9.97	3390420	20.0
unknown10.24	10.24	2.6	ng	444877	3	9.97	3390420	20.0
unknown10.30	10.30	2.5	ng	414948	3	9.97	3390420	20.0
1-Trimethylsilylp...	10.39	3.8	ng	645074	3	9.97	3390420	20.0
Pentadecane, 2,6,...	10.63	7.0	ng	1184950	3	9.97	3390420	20.0
Tridecane, 5-propyl-	10.90	29.7	ng	4013290	4	11.47	2700530	20.0
Hexadecane, 2,6,1...	11.37	22.7	ng	3067420	4	11.47	2700530	20.0
1-Octadecanol	13.96	18.6	ng	1761710	5	14.12	1897460	20.0
Supraene	14.99	13.1	ng	1461420	6	15.62	2225420	20.0
Benzo[e]pyrene	15.49	5.5	ng	609569	6	15.62	2225420	20.0