

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121618\
 Data File : BF111526.D
 Acq On : 16 Dec 2018 23:42
 Operator : JU/SJ
 Sample : J6377-04
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 P001-SD014-0612-01

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-BF121418.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	2.887	128	136	142	rVB	44501	104039	1.24%	0.212%
2	3.822	286	295	301	rBV	366890	524579	6.26%	1.069%
3	4.422	391	397	402	rVB2	58492	84139	1.00%	0.172%
4	4.998	492	495	501	rBV	157541	200136	2.39%	0.408%
5	5.081	505	509	513	rBV	242908	241488	2.88%	0.492%
6	5.269	537	541	544	rBV	465001	431262	5.15%	0.879%
7	5.387	555	561	564	rBV	6192284	8374310	100.00%	17.070%
8	5.422	564	567	570	rVB	3982011	3528504	42.13%	7.192%
9	5.634	600	603	607	rBV	710397	642866	7.68%	1.310%
10	6.316	717	719	722	rBV	178862	149779	1.79%	0.305%
11	6.392	728	732	735	rBV2	159105	139730	1.67%	0.285%
12	6.434	735	739	744	rBV2	6038594	6126361	73.16%	12.488%
13	6.563	757	761	764	rVV	3816304	3500492	41.80%	7.135%
14	6.616	767	770	776	rVB2	498151	522959	6.24%	1.066%
15	6.787	796	799	801	rBV	866007	809784	9.67%	1.651%
16	6.804	801	802	807	rVB	743405	528364	6.31%	1.077%
17	6.851	807	810	818	rVB2	157196	173509	2.07%	0.354%
18	6.945	822	826	829	rBV	3190799	2765049	33.02%	5.636%
19	7.069	844	847	850	rVB2	102238	93759	1.12%	0.191%
20	7.110	850	854	856	rBV2	116673	105019	1.25%	0.214%
21	7.351	886	895	898	rBV	2213301	2334549	27.88%	4.759%
22	7.392	900	902	906	rVV	218579	171499	2.05%	0.350%
23	7.563	926	931	933	rBV	506611	429381	5.13%	0.875%
24	7.586	933	935	940	rVB	93189	93997	1.12%	0.192%
25	8.069	1012	1017	1020	rBV	1097429	1036076	12.37%	2.112%
26	8.610	1106	1109	1112	rBV	129800	118764	1.42%	0.242%
27	8.669	1115	1119	1122	rBV	166581	130847	1.56%	0.267%
28	8.780	1132	1138	1142	rVB2	105578	110480	1.32%	0.225%
29	9.145	1196	1200	1203	rBV	4167236	3692837	44.10%	7.527%
30	9.233	1211	1215	1220	rBV2	218073	207251	2.47%	0.422%
31	9.539	1264	1267	1268	rBV	186028	168027	2.01%	0.342%
32	9.763	1303	1305	1309	rVB	166162	123710	1.48%	0.252%
33	9.822	1309	1315	1319	rBV	988240	915959	10.94%	1.867%
34	10.263	1387	1390	1393	rVB	206463	169559	2.02%	0.346%

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ClientSampleId :
P001-SD014-0612-01

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	10.310	1393	1398	1402	rBV2	76601	84989	1.01%	0.173%
36	10.480	1421	1427	1431	rBV	92807	118400	1.41%	0.241%
37	10.610	1445	1449	1462	rBV	1703488	1592220	19.01%	3.245%
38	10.745	1467	1472	1480	rVB2	142266	243606	2.91%	0.497%
39	11.180	1544	1546	1549	rBV	103399	107403	1.28%	0.219%
40	11.210	1549	1551	1555	rVV2	90888	99087	1.18%	0.202%
41	11.304	1564	1567	1570	rBV	1051338	952714	11.38%	1.942%
42	11.839	1655	1658	1661	rBV	125033	130976	1.56%	0.267%
43	12.298	1731	1736	1741	rBV3	59437	111417	1.33%	0.227%
44	12.892	1833	1837	1847	rBV	3838740	3650960	43.60%	7.442%
45	13.080	1866	1869	1871	rVB	106164	91251	1.09%	0.186%
46	13.268	1898	1901	1904	rVB2	131882	124717	1.49%	0.254%
47	13.374	1916	1919	1925	rVB6	63509	95818	1.14%	0.195%
48	13.474	1933	1936	1942	rVB2	114160	148223	1.77%	0.302%
49	13.586	1952	1955	1960	rBV5	119490	111291	1.33%	0.227%
50	13.774	1984	1987	1988	rBV2	108349	102945	1.23%	0.210%
51	13.798	1988	1991	2002	rVB3	206957	375565	4.48%	0.766%
52	13.945	2012	2016	2019	rBV	1008483	1000218	11.94%	2.039%
53	13.980	2019	2022	2027	rVB4	127003	126327	1.51%	0.257%
54	14.421	2094	2097	2102	rVV2	121008	175670	2.10%	0.358%
55	14.515	2110	2113	2118	rBV2	84815	129123	1.54%	0.263%
56	15.051	2201	2204	2206	rBV	100177	104634	1.25%	0.213%
57	15.380	2256	2260	2264	rBV	569105	632796	7.56%	1.290%

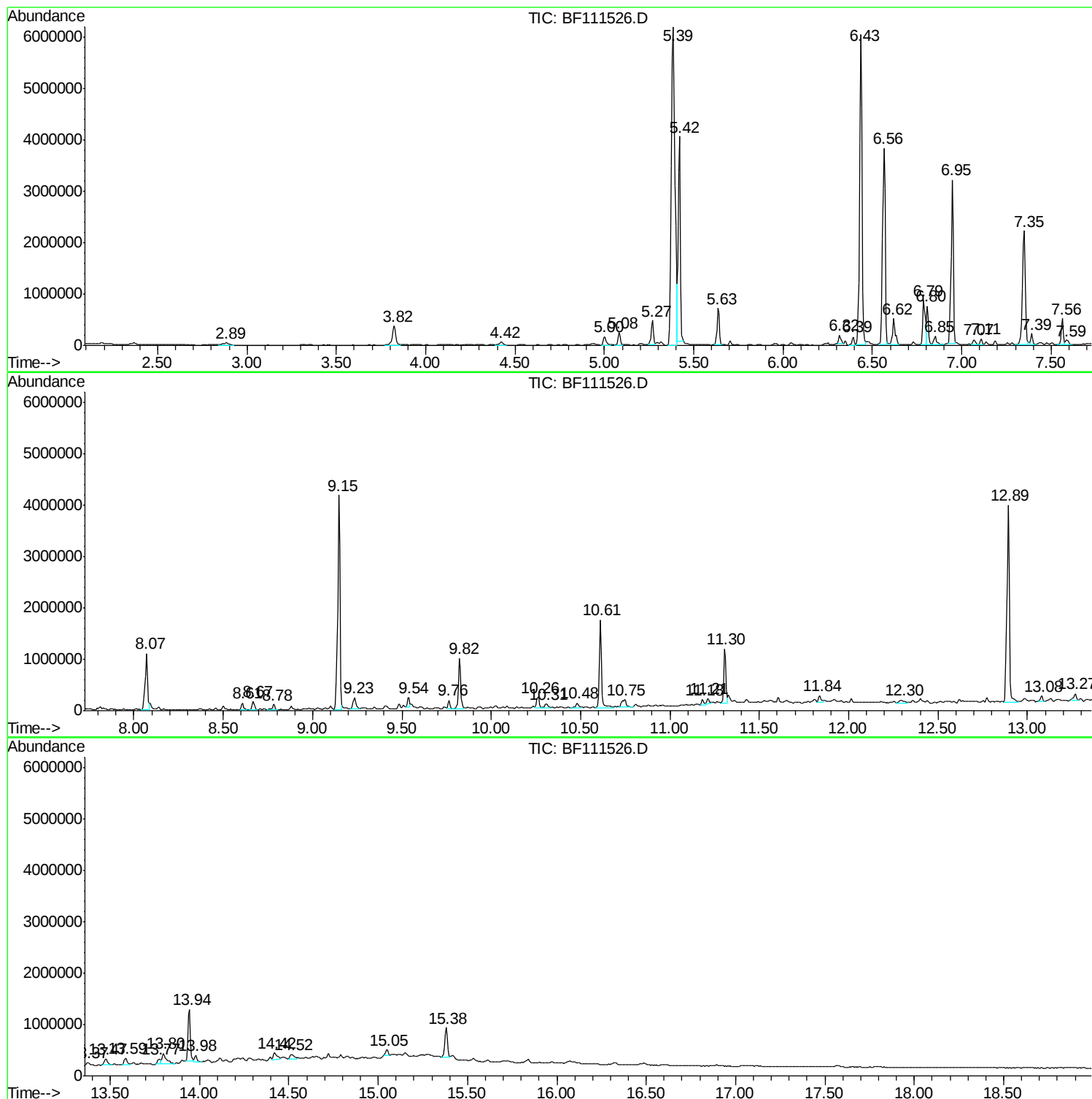
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF121418.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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Operator : JU/SJ
Sample : J6377-04
Misc :
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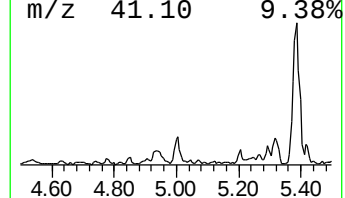
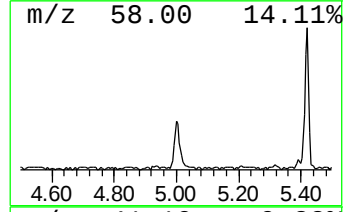
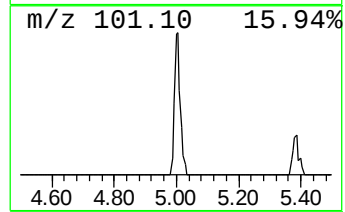
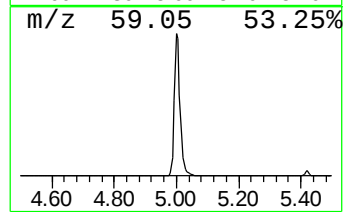
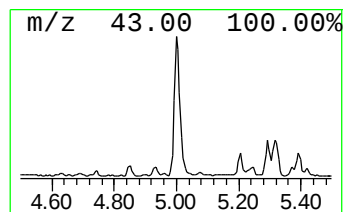
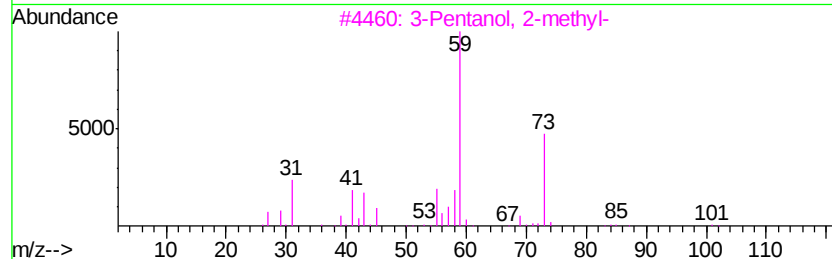
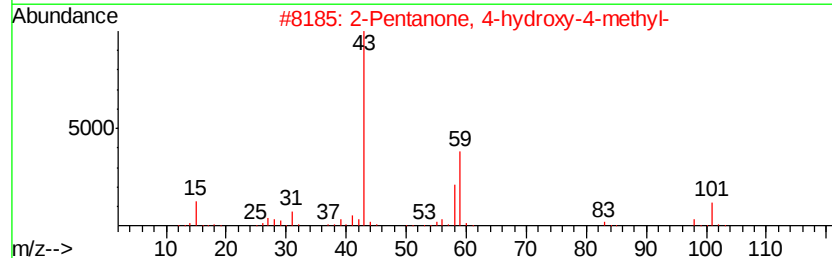
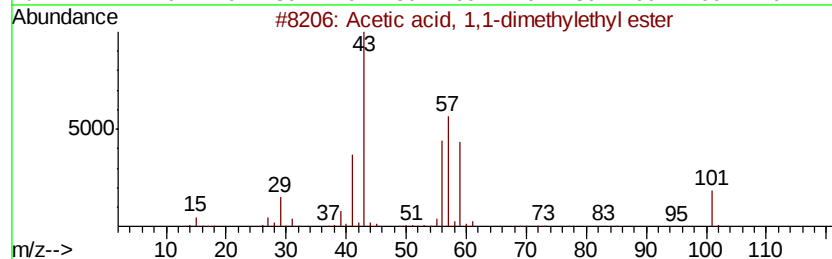
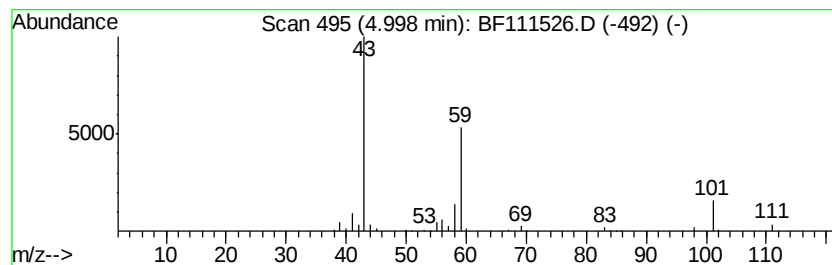
Instrument :
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ClientSampleId :
P001-SD014-0612-01

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF121418.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 Acetic acid, 1,1-dimethylet... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.		
5.00	4.94 ng	200136	1,4-Dichlorobenzene-d4	6.79		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	50	
2	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50	
3	3-Pentanol, 2-methyl-	102	C6H14O	000565-67-3	28	
4	2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	25	
5	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	25	



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Instrument :
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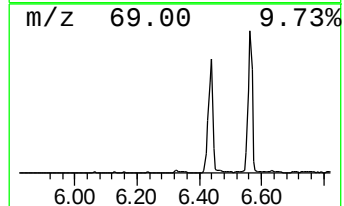
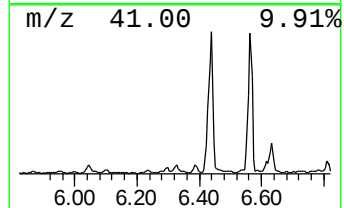
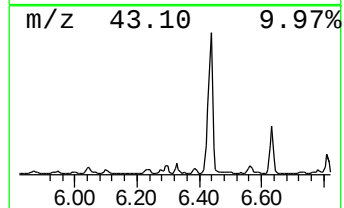
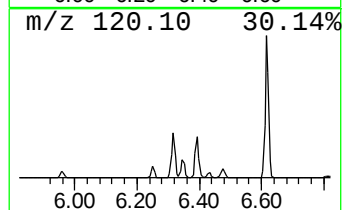
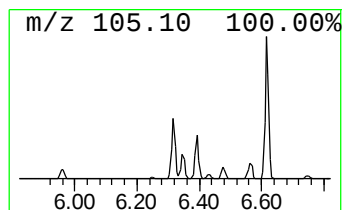
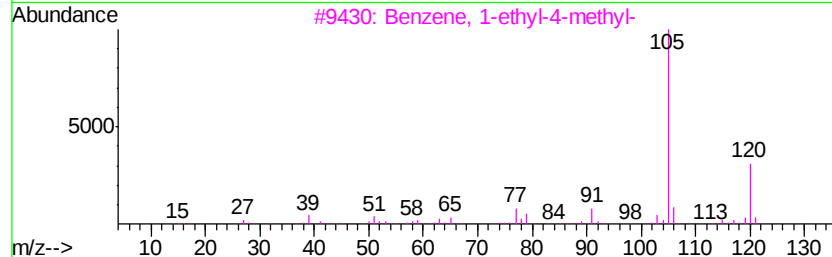
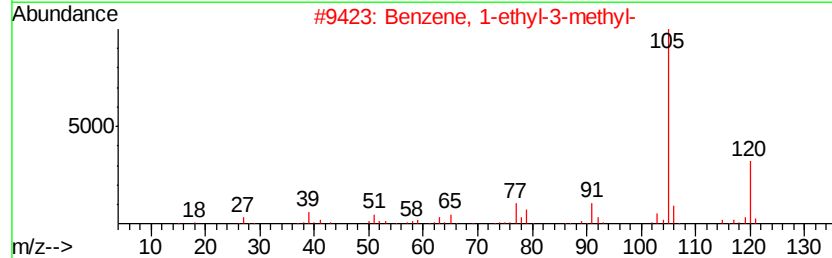
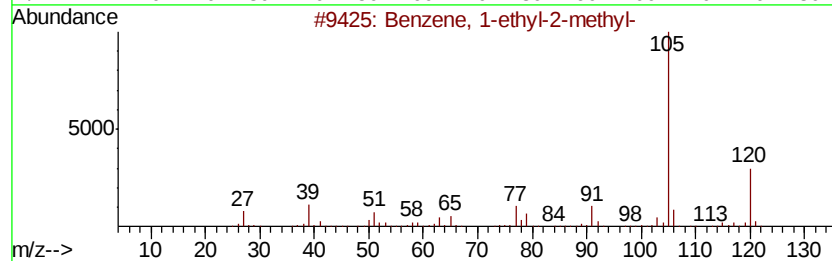
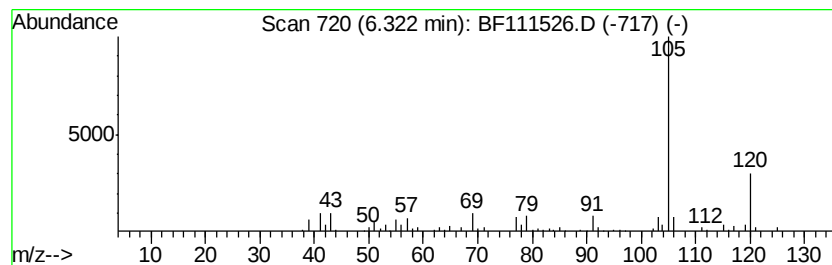
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF121418.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Benzene, 1-ethyl-2-methyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.32	3.70 ng	149779	1,4-Dichlorobenzene-d4	6.79

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
2		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	95
3		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
4		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
5		Mesitylene	120	C9H12	000108-67-8	91



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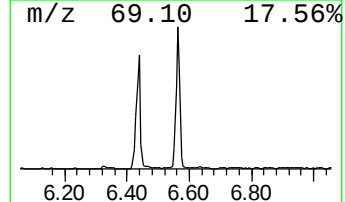
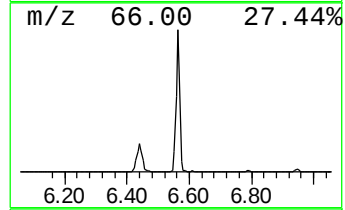
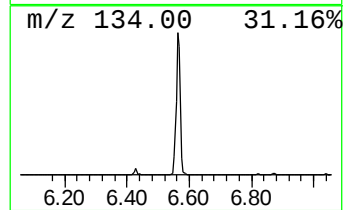
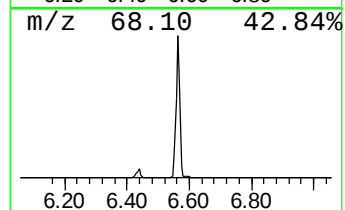
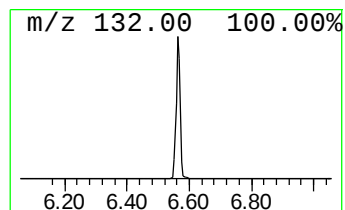
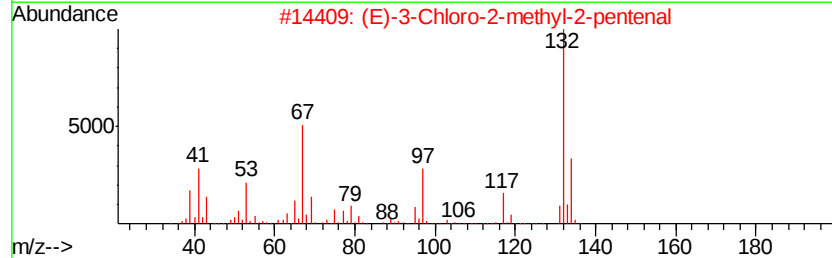
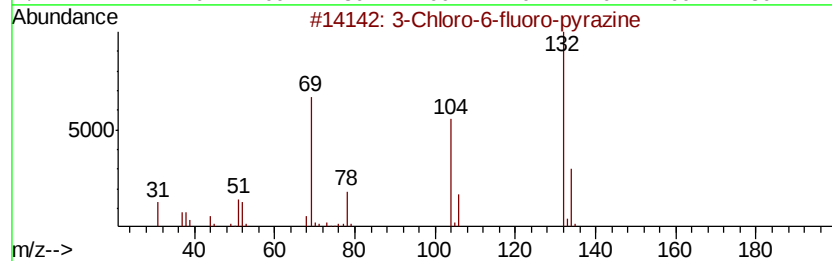
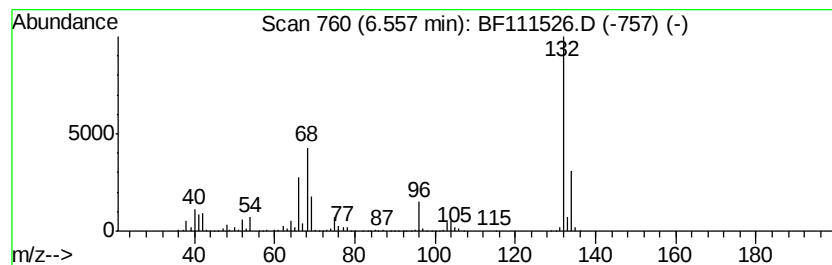
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 unknown6.56 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.56	86.46 ng	3500490	1,4-Dichlorobenzene-d4	6.79

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	35
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
3		Tranvlcyproline-propionyl	189	C12H15NO	1000123-86-3	10
4		1,3,5-Trifluorobenzene	132	C6H3F3	000372-38-3	9
5		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9



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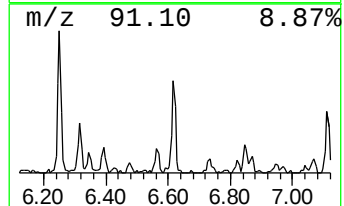
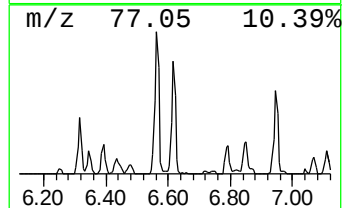
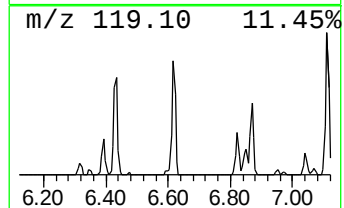
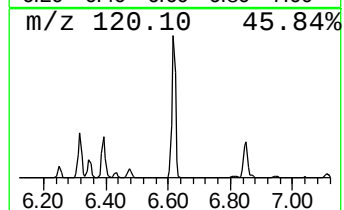
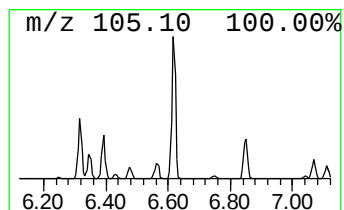
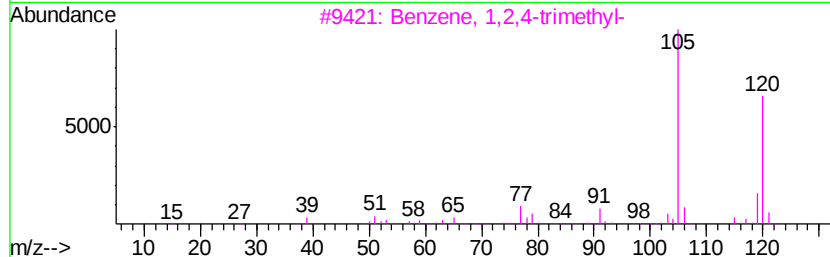
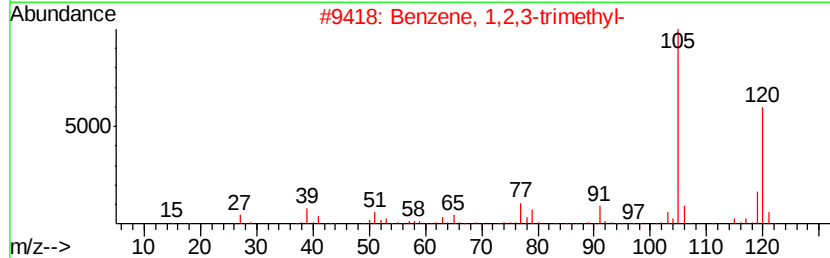
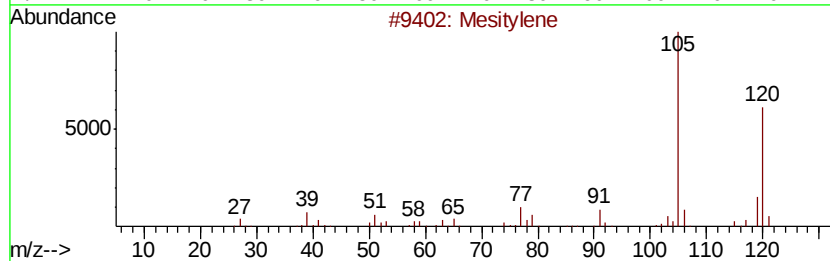
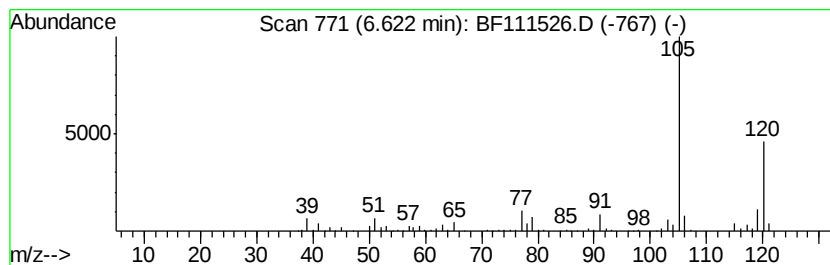
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Mesitylene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.62	12.92 ng	522959	1,4-Dichlorobenzene-d4	6.79

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Mesitylene	120	C9H12	000108-67-8	97
2		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	97
3		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94
4		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	93
5		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	90



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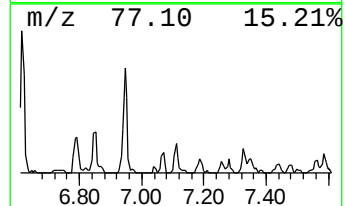
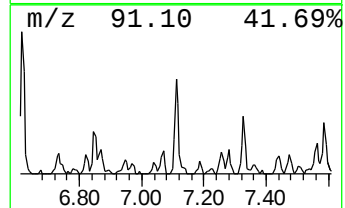
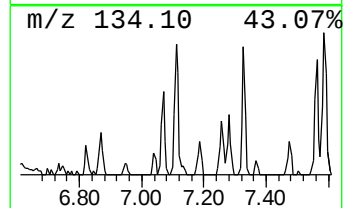
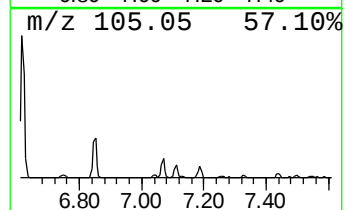
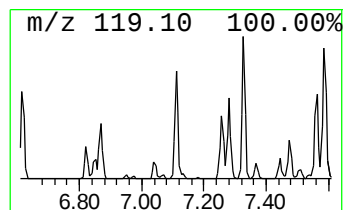
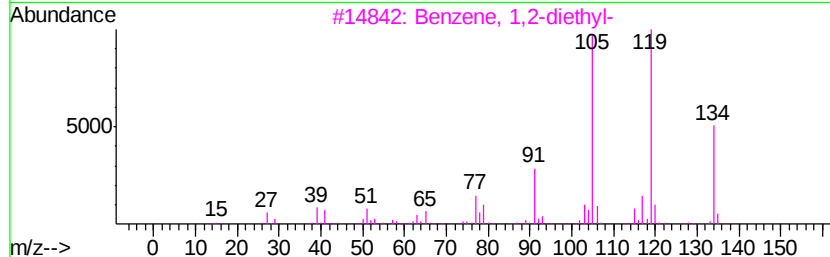
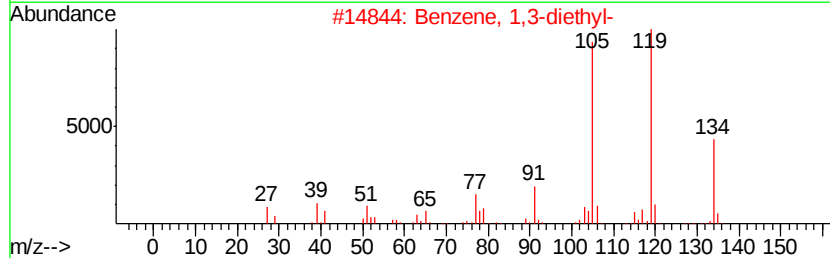
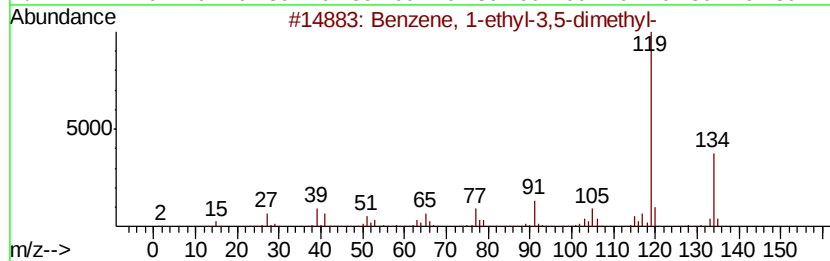
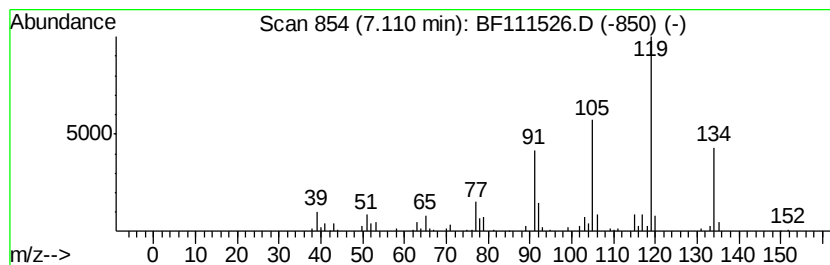
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Benzene, 1-ethyl-3,5-dimethyl- Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.11	2.59 ng	105019	1,4-Dichlorobenzene-d4	6.79

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	96
2		Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	94
3		Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	92
4		1,3,8-p-Menthatriene	134	C10H14	018368-95-1	91
5		Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121618\
Data File : BF111526.D
Acq On : 16 Dec 2018 23:42
Operator : JU/SJ
Sample : J6377-04
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
P001-SD014-0612-01

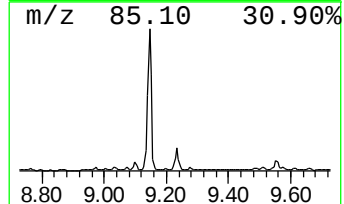
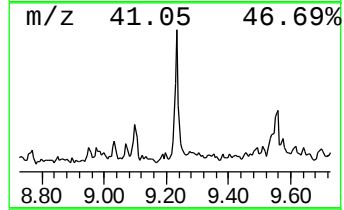
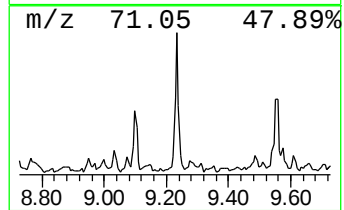
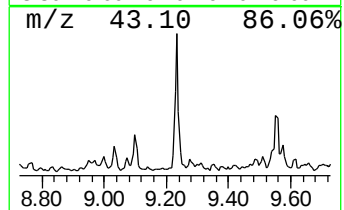
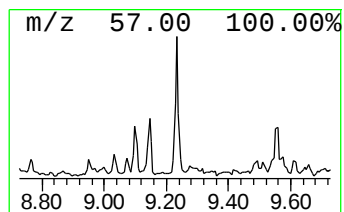
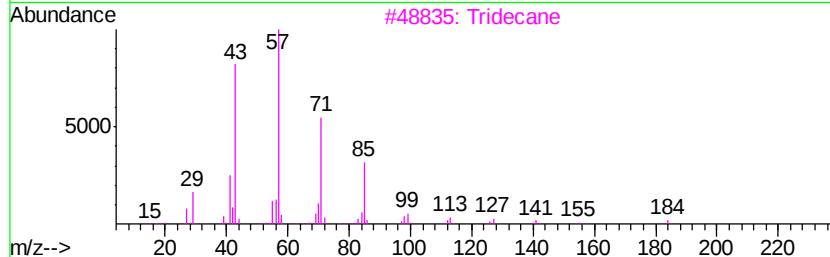
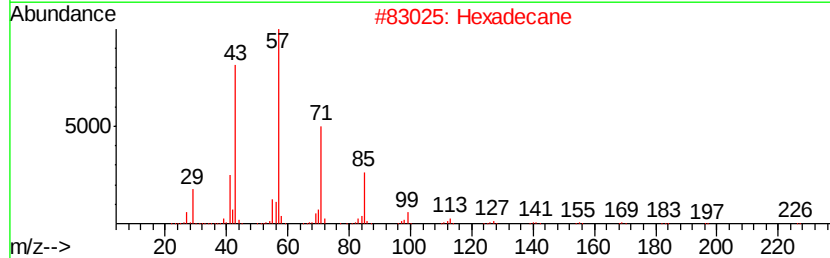
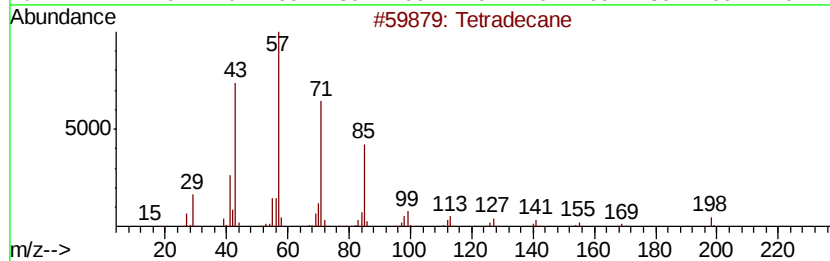
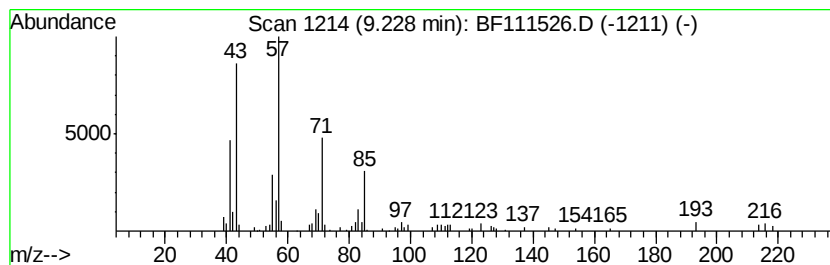
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Tetradecane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.23	4.53 ng	207251	Acenaphthene-d10	9.82

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetradecane	198	C14H30	000629-59-4	87
2		Hexadecane	226	C16H34	000544-76-3	83
3		Tridecane	184	C13H28	000629-50-5	80
4		Dodecane	170	C12H26	000112-40-3	80
5		Undecane	156	C11H24	001120-21-4	80



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121618\
Data File : BF111526.D
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Operator : JU/SJ
Sample : J6377-04
Misc :
ALS Vial : 23 Sample Multiplier: 1

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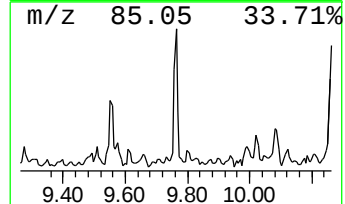
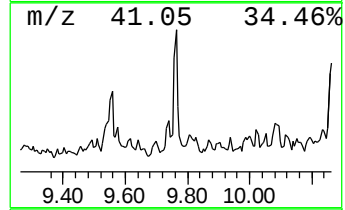
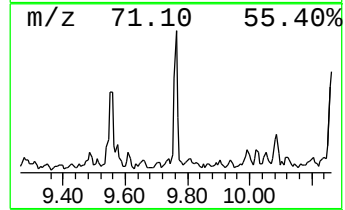
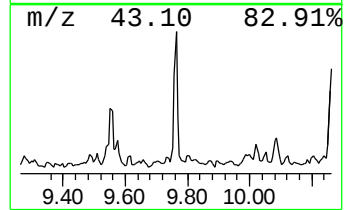
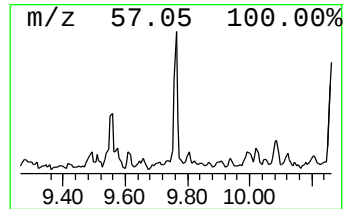
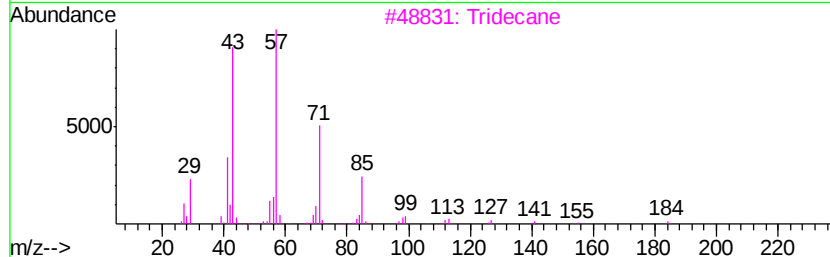
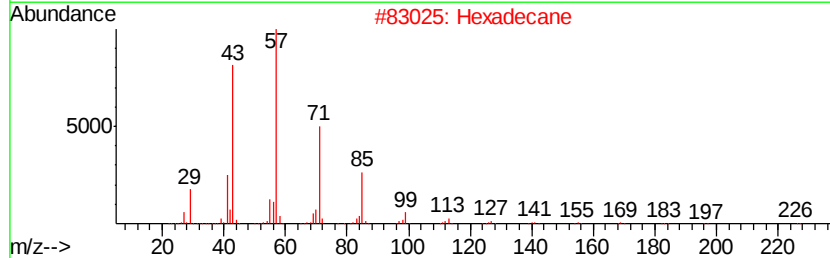
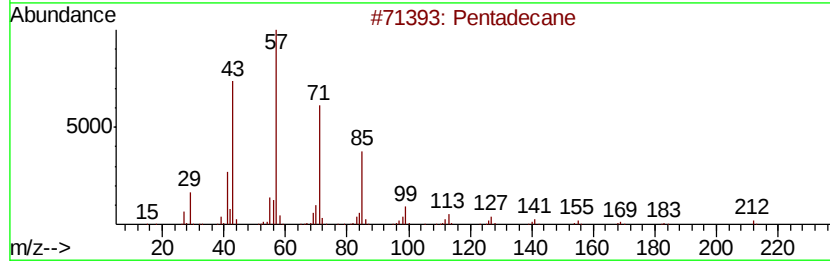
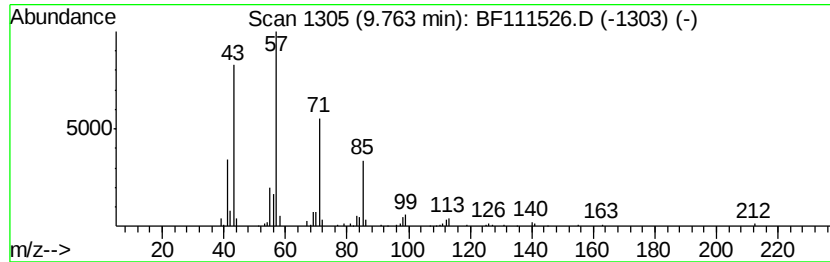
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Pentadecane Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.76	2.70 ng	123710	Acenaphthene-d10	9.82

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentadecane	212	C15H32	000629-62-9	91
2			Hexadecane	226	C16H34	000544-76-3	86
3			Tridecane	184	C13H28	000629-50-5	78
4			Tridecane, 4,8-dimethyl-	212	C15H32	055030-62-1	72
5			Decane, 3-methyl-	156	C11H24	013151-34-3	72



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121618\
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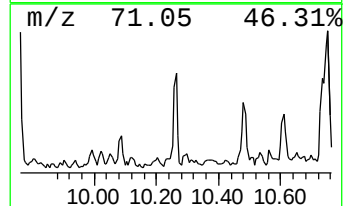
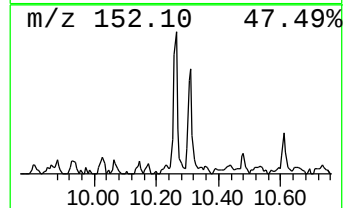
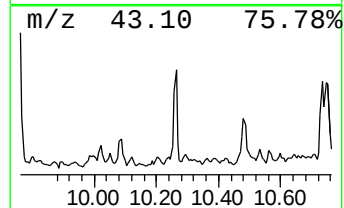
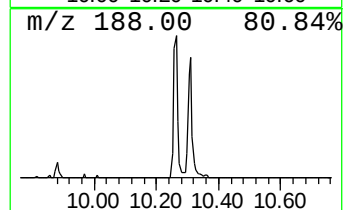
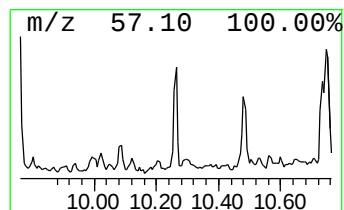
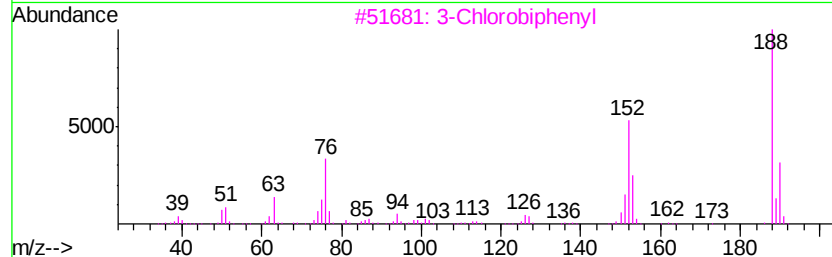
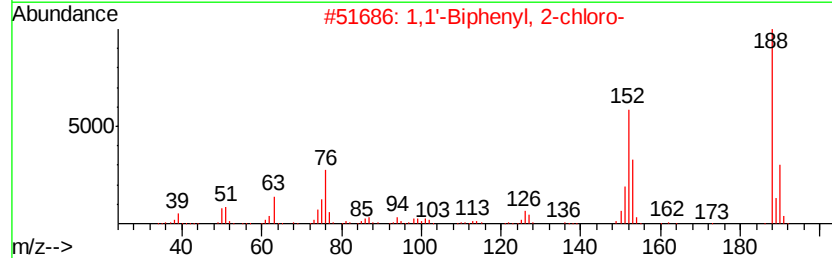
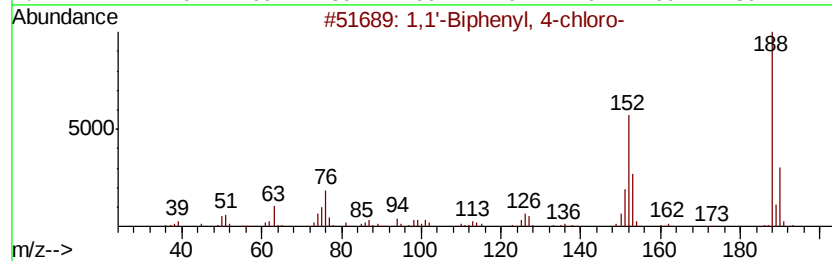
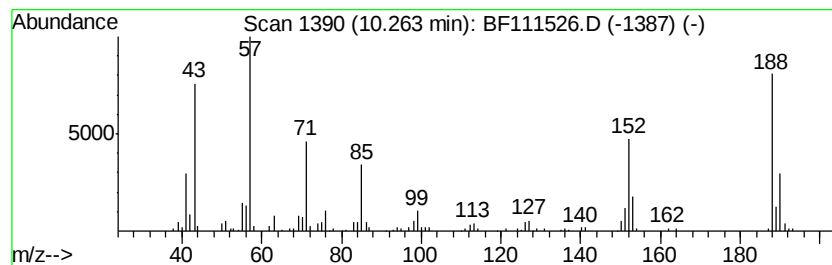
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 1,1'-Biphenyl, 4-chloro- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD		R.T.
10.26	3.70 ng	169559	Acenaphthene-d10		9.82
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 4-chloro-	188	C12H9Cl	002051-62-9	92
2	1,1'-Biphenyl, 2-chloro-	188	C12H9Cl	002051-60-7	78
3	3-Chlorobiphenyl	188	C12H9Cl	002051-61-8	64
4	3,5-Dimethyldodecane	198	C14H30	107770-99-0	11
5	Hexane, 2,3,4-trimethyl-	128	C9H20	000921-47-1	11



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121618\
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Sample : J6377-04
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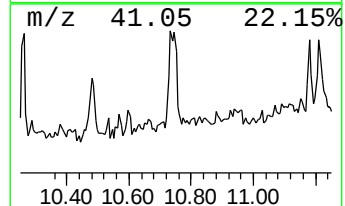
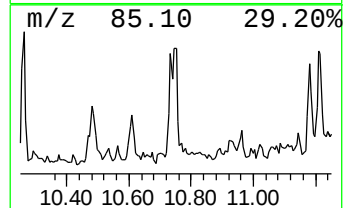
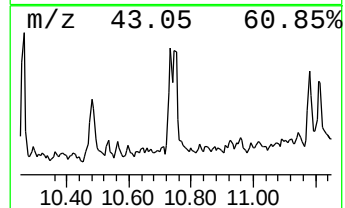
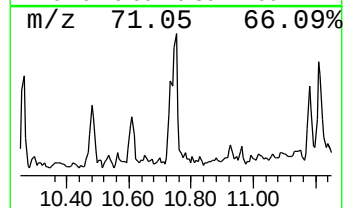
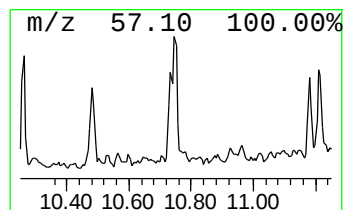
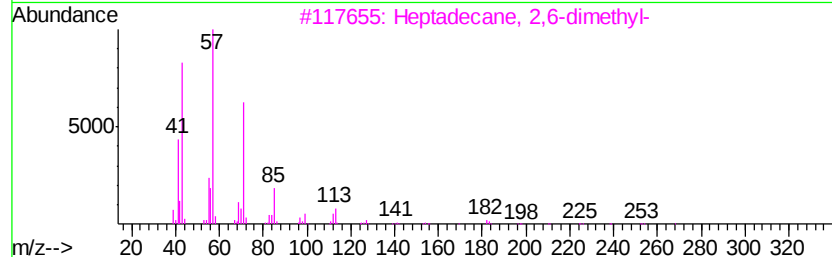
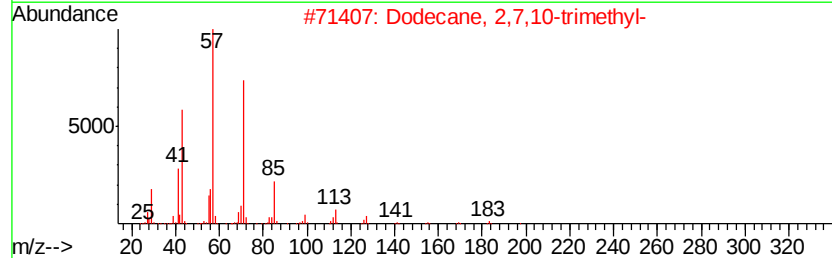
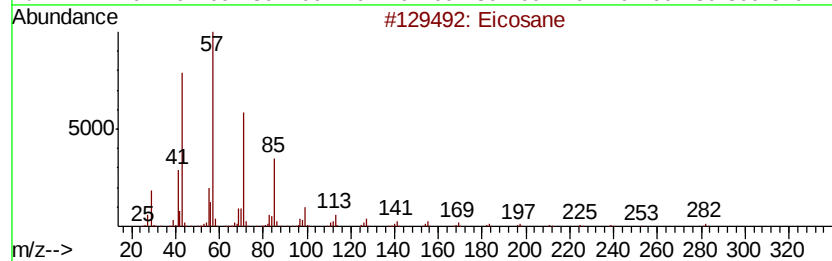
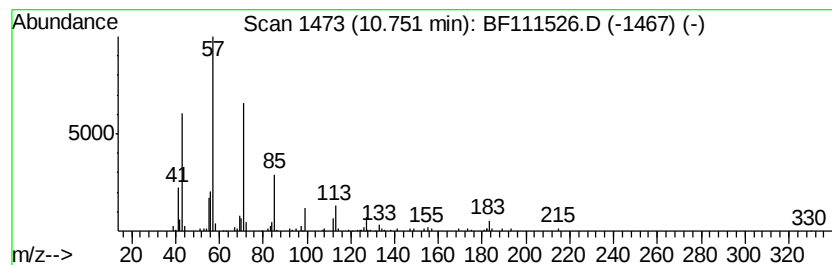
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Eicosane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD		R.T.	
10.75	5.11 ng	243606	Phenanthrene-d10		11.30	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane		282	C20H42	000112-95-8	72
2	Dodecane, 2,7,10-trimethyl-		212	C15H32	074645-98-0	64
3	Heptadecane, 2,6-dimethyl-		268	C19H40	054105-67-8	64
4	Pentadecane, 2,6,10,14-tetramethyl-		268	C19H40	001921-70-6	64
5	Bacchotricuneatin c		342	C20H22O5	066563-30-2	60



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121618\
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Sample : J6377-04
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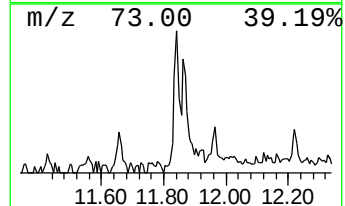
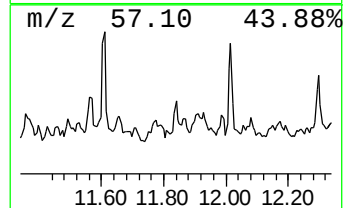
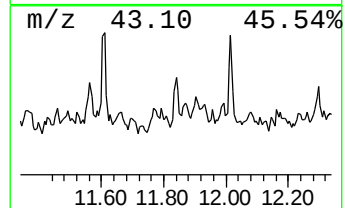
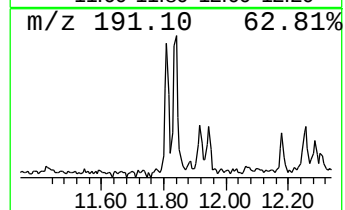
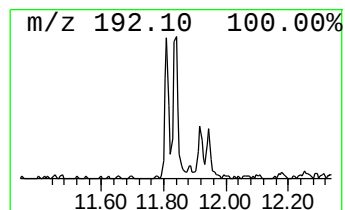
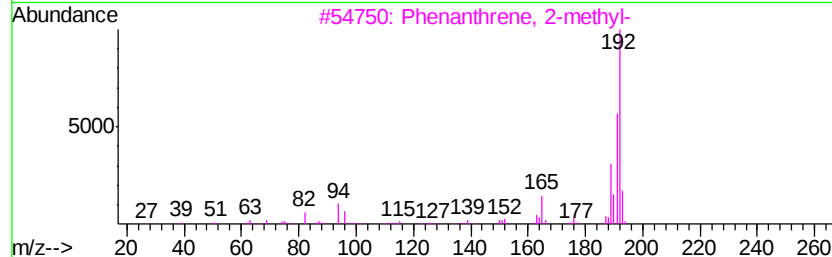
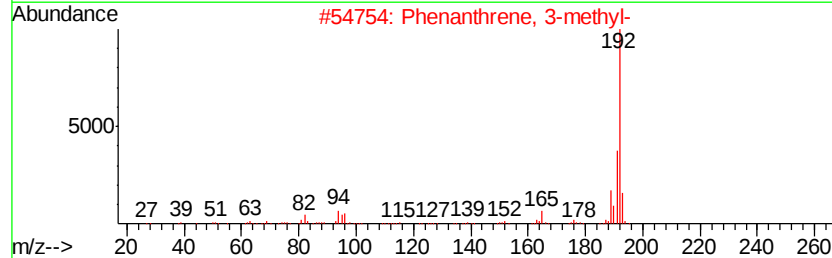
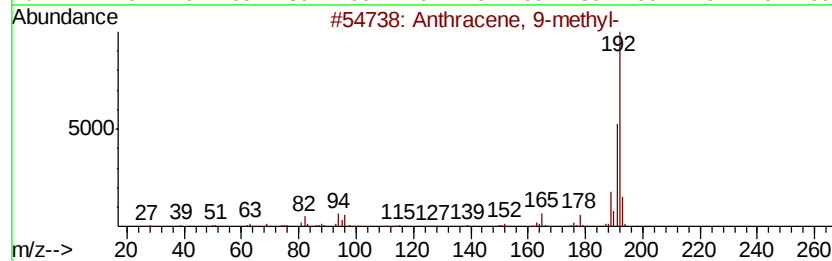
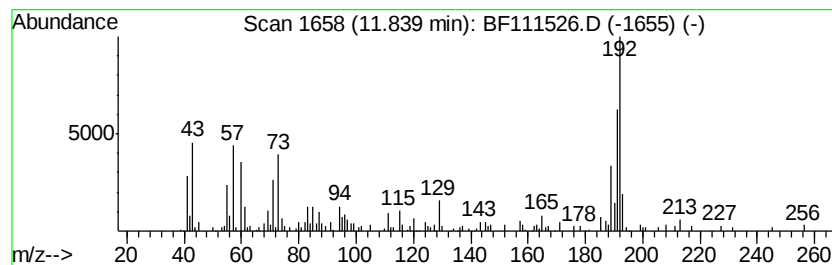
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 Anthracene, 9-methyl- Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.84	2.75 ng	130976	Phenanthrene-d10	11.30	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Anthracene, 9-methyl-	192	C15H12	000779-02-2	93
2	Phenanthrene, 3-methyl-	192	C15H12	000832-71-3	90
3	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	83
4	1H-Cyclopropa[1]phenanthrene,1a,...	192	C15H12	000949-41-7	78
5	Anthracene, 2-methyl-	192	C15H12	000613-12-7	70



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121618\
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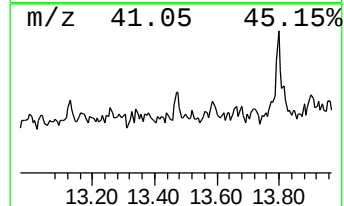
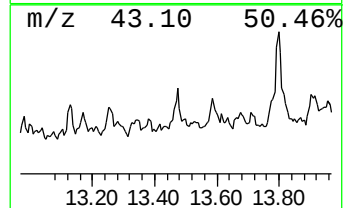
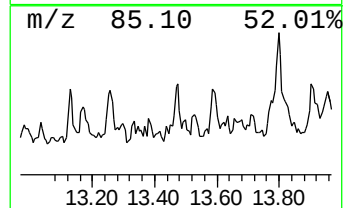
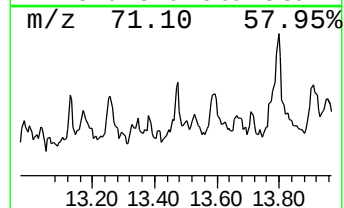
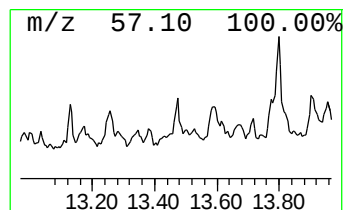
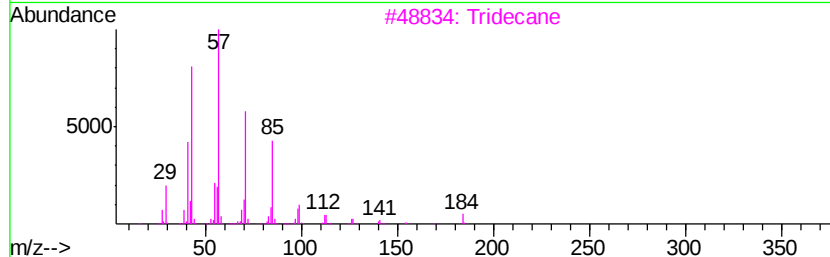
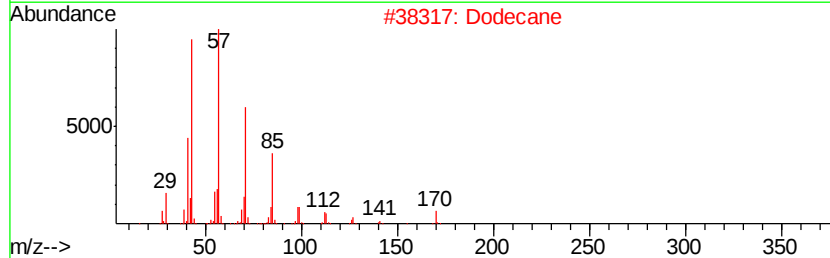
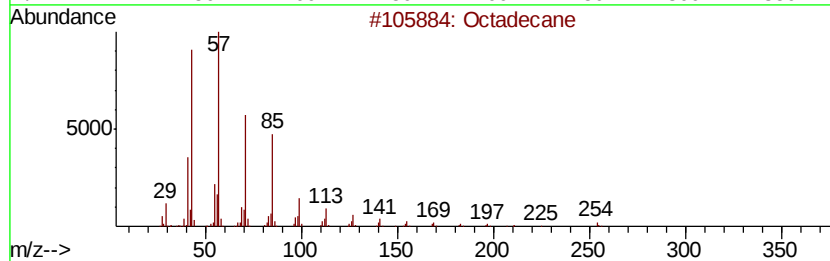
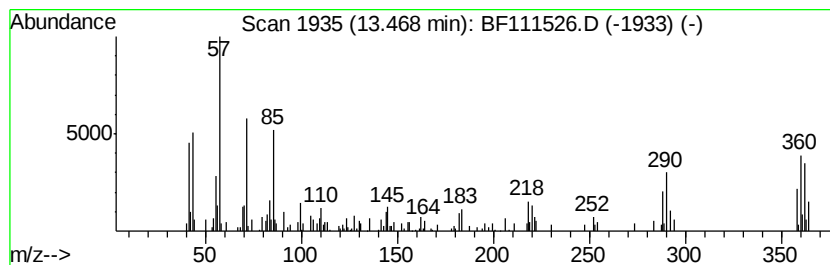
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TIC Library : C:\DATABASE\NIST11.L
 TIC Integration Parameters: LSCINT.P

 Peak Number 17 unknown13.47 Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.47	2.96 ng	148223	Chrysene-d12	13.94

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecane	254	C18H38	000593-45-3	45
2		Dodecane	170	C12H26	000112-40-3	40
3		Tridecane	184	C13H28	000629-50-5	25
4		Undecane	156	C11H24	001120-21-4	15
5		Benzoic acid, 2-(1-methyl-4-nitr...	290	C12H10N4O5	1000295-63-7	14



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121618\
Data File : BF111526.D
Acq On : 16 Dec 2018 23:42
Operator : JU/SJ
Sample : J6377-04
Misc :
ALS Vial : 23 Sample Multiplier: 1

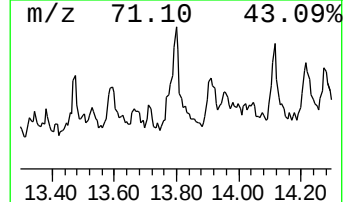
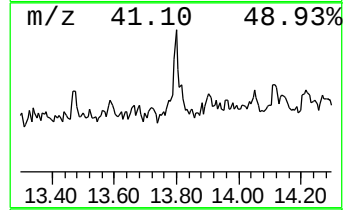
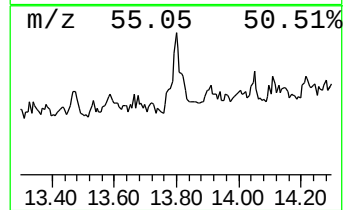
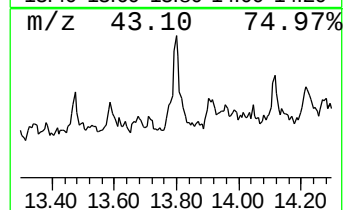
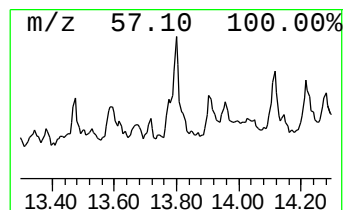
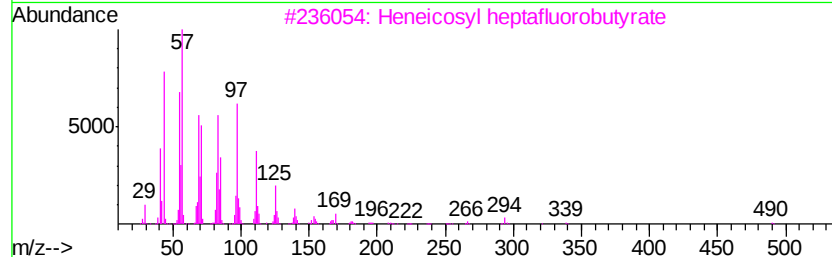
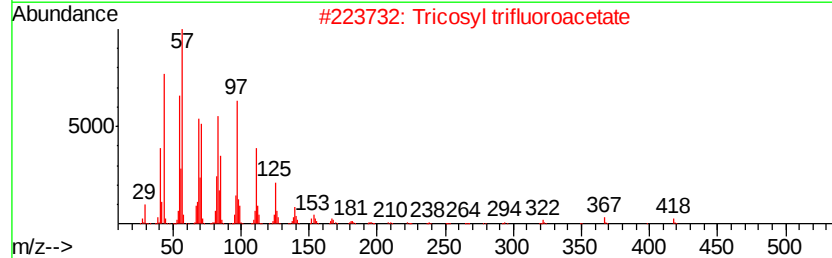
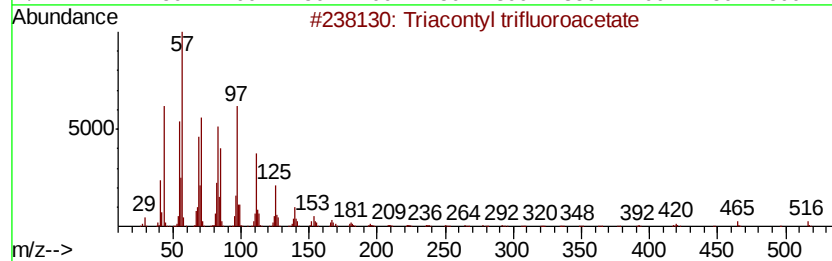
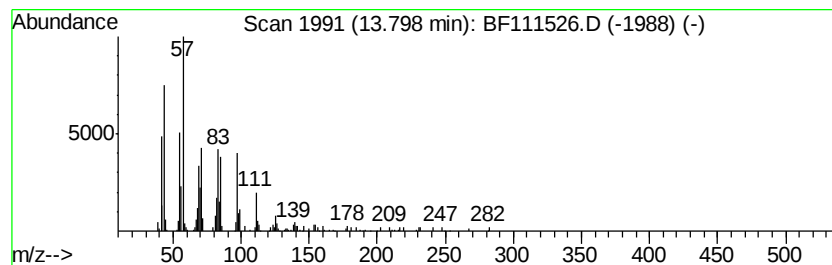
Instrument :
BNA_F
ClientSampleId :
P001-SD014-0612-01

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF121418.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 18 Triacontyl trifluoroacetate Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD		R.T.	
13.80	7.51 ng	375565	Chrysene-d12		13.94	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Triacontyl trifluoroacetate	534	C32H61F3O2	1000351-75-7	90
2		Tricosyl trifluoroacetate	436	C25H47F3O2	1000351-75-1	86
3		Heneicosyl heptafluorobutvrate	508	C25H43F7O2	1000351-83-8	83
4		Docosyl pentafluoropropionate	472	C25H45F5O2	1000351-80-9	83
5		Oxalic acid, propyl tetradecyl e...	328	C19H36O4	1000309-26-7	74



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121618\
Data File : BF111526.D
Acq On : 16 Dec 2018 23:42
Operator : JU/SJ
Sample : J6377-04
Misc :
ALS Vial : 23 Sample Multiplier: 1

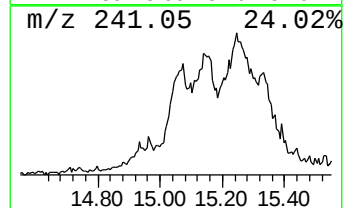
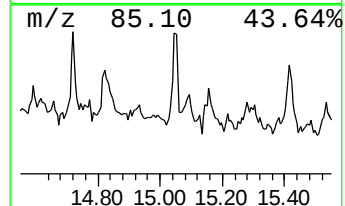
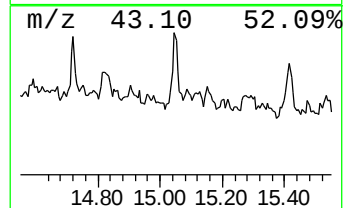
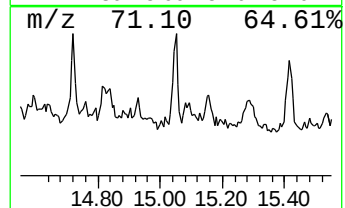
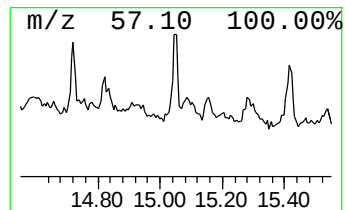
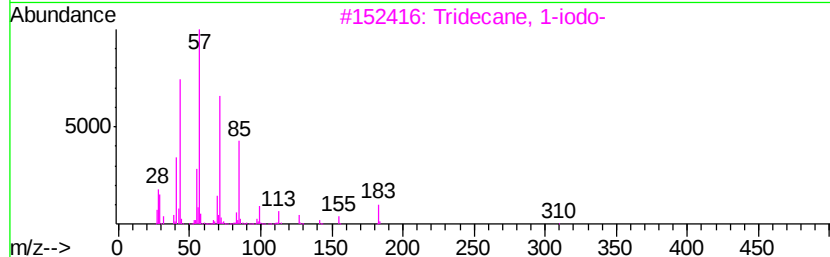
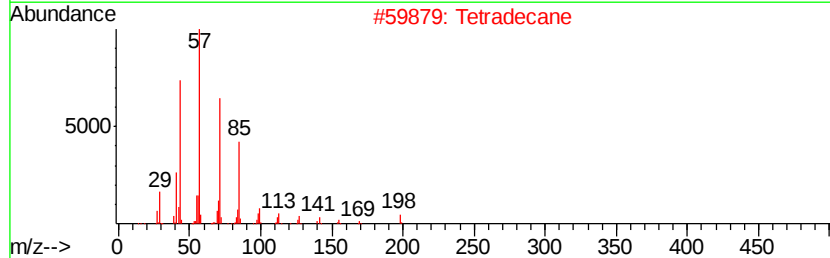
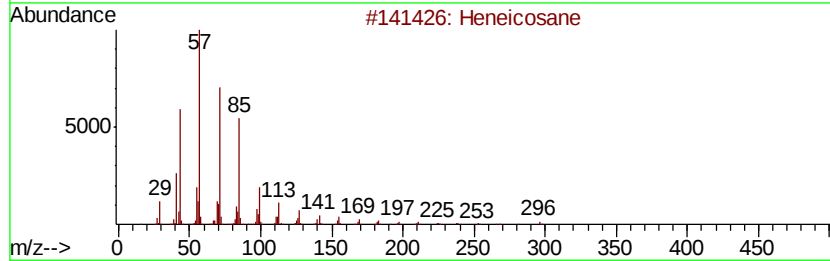
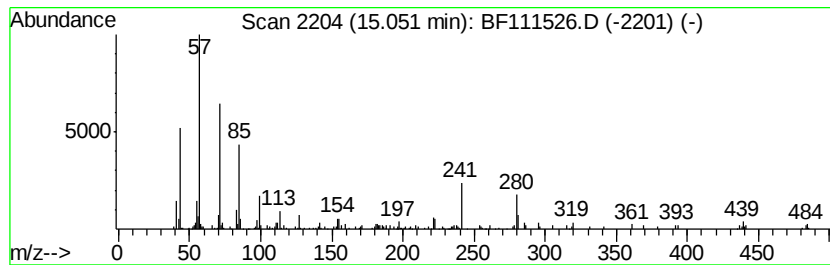
Instrument :
BNA_F
ClientSampleId :
P001-SD014-0612-01

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF121418.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 20 Heneicosane Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.05	3.31 ng	104634	Perylene-d12	15.38
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Heneicosane	296 C21H44	000629-94-7	60
2	Tetradecane	198 C14H30	000629-59-4	53
3	Tridecane, 1-iodo-	310 C13H27I	035599-77-0	49
4	Dodecane, 2-methyl-	184 C13H28	001560-97-0	49
5	Octadecane	254 C18H38	000593-45-3	49



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF121618\
 Data File : BF111526.D
 Acq On : 16 Dec 2018 23:42
 Operator : JU/SJ
 Sample : J6377-04
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 P001-SD014-0612-01

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF121418.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
Acetic acid, 1,1-...	5.00	4.9	ng	200136	1	6.79	809784	20.0
Benzene, 1-ethyl-...	6.32	3.7	ng	149779	1	6.79	809784	20.0
unknown6.56	6.56	86.5	ng	3500490	1	6.79	809784	20.0
Mesitylene	6.62	12.9	ng	522959	1	6.79	809784	20.0
Benzene, 1-ethyl-...	7.11	2.6	ng	105019	1	6.79	809784	20.0
Tetradecane	9.23	4.5	ng	207251	3	9.82	915959	20.0
Pentadecane	9.76	2.7	ng	123710	3	9.82	915959	20.0
1,1'-Biphenyl, 4-...	10.26	3.7	ng	169559	3	9.82	915959	20.0
Eicosane	10.75	5.1	ng	243606	4	11.30	952714	20.0
Anthracene, 9-met...	11.84	2.8	ng	130976	4	11.30	952714	20.0
unknown13.47	13.47	3.0	ng	148223	5	13.94	1000220	20.0
Triacontyl triflu...	13.80	7.5	ng	375565	5	13.94	1000220	20.0
Heneicosane	15.05	3.3	ng	104634	6	15.38	632796	20.0