

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF021819\
 Data File : BF112603.D
 Acq On : 18 Feb 2019 16:13
 Operator : JU/SJ
 Sample : K1518-15
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CABLE-BRIDGE-FOUNDATION-D-8

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-BF012519.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	5.028	496	500	513	rBV	39683	57836	2.51%	0.313%
2	5.457	569	573	597	rBV	865583	1081876	46.93%	5.862%
3	5.657	604	607	616	rBV2	26290	36275	1.57%	0.197%
4	6.475	742	746	764	rBV	1012723	1238839	53.73%	6.712%
5	6.598	764	767	773	rVV	1382075	1278574	55.46%	6.927%
6	6.657	773	777	781	rVB2	44038	51745	2.24%	0.280%
7	6.822	802	805	812	rBV	495328	421603	18.29%	2.284%
8	6.981	828	832	847	rBV	1056805	989467	42.92%	5.361%
9	7.087	847	850	857	rVV2	23616	27240	1.18%	0.148%
10	7.387	897	901	909	rBV	769059	758478	32.90%	4.109%
11	7.681	945	951	959	rBV	64204	80271	3.48%	0.435%
12	8.104	1020	1023	1030	rVB	590593	525119	22.78%	2.845%
13	8.322	1053	1060	1066	rBV	54398	66013	2.86%	0.358%
14	8.381	1066	1070	1079	rVV	63614	82088	3.56%	0.445%
15	8.522	1091	1094	1096	rBV	38919	32727	1.42%	0.177%
16	8.692	1118	1123	1126	rBV	42022	40264	1.75%	0.218%
17	8.816	1141	1144	1148	rBV3	21683	27441	1.19%	0.149%
18	8.986	1169	1173	1176	rBV3	26629	33907	1.47%	0.184%
19	9.122	1194	1196	1202	rVB2	47624	32465	1.41%	0.176%
20	9.175	1202	1205	1214	rBV	1707322	1528939	66.32%	8.284%
21	9.257	1215	1219	1222	rBV	49821	50640	2.20%	0.274%
22	9.575	1269	1273	1282	rBV	156513	183719	7.97%	0.995%
23	9.728	1295	1299	1301	rBV	51595	54577	2.37%	0.296%
24	9.863	1318	1322	1326	rVB	691256	638097	27.68%	3.457%
25	10.281	1387	1393	1396	rBV4	34488	41776	1.81%	0.226%
26	10.657	1453	1457	1470	rBV	1063410	1036218	44.95%	5.614%
27	10.769	1471	1476	1478	rVV2	25510	30737	1.33%	0.167%
28	11.351	1571	1575	1578	rBV	751448	654606	28.39%	3.547%
29	11.592	1611	1616	1620	rVB2	20416	26965	1.17%	0.146%
30	11.869	1659	1663	1671	rBV3	229204	277156	12.02%	1.502%
31	12.192	1715	1718	1725	rVB6	24937	29204	1.27%	0.158%
32	12.569	1778	1782	1786	rBV	268102	231884	10.06%	1.256%
33	12.651	1793	1796	1803	rBV3	85980	115906	5.03%	0.628%
34	12.798	1818	1821	1827	rVB	252708	228458	9.91%	1.238%

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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	12.927	1839	1843	1847	rBV	1604576	1398563	60.66%	7.577%
36	13.104	1871	1873	1876	rBV3	21434	26313	1.14%	0.143%
37	13.492	1936	1939	1942	rBV4	27201	30346	1.32%	0.164%
38	13.645	1963	1965	1970	rVB	35444	39465	1.71%	0.214%
39	13.751	1981	1983	1985	rBV2	27997	29659	1.29%	0.161%
40	13.816	1989	1994	1999	rVV3	98795	164314	7.13%	0.890%
41	13.857	1999	2001	2007	rVB2	38973	46944	2.04%	0.254%
42	13.957	2014	2018	2021	rBV	2616183	2305467	100.00%	12.491%
43	13.998	2021	2025	2028	rBV	604663	604307	26.21%	3.274%
44	14.133	2045	2048	2050	rBV	76495	64197	2.78%	0.348%
45	14.439	2097	2100	2107	rVB3	125895	127143	5.51%	0.689%
46	14.739	2148	2151	2155	rVB	131286	128647	5.58%	0.697%
47	15.045	2200	2203	2205	rBV	135086	167968	7.29%	0.910%
48	15.068	2205	2207	2214	rVB3	188569	253153	10.98%	1.372%
49	15.445	2268	2271	2273	rBV	82390	103274	4.48%	0.560%
50	15.474	2273	2276	2282	rVB2	361013	464501	20.15%	2.517%
51	15.863	2336	2342	2351	rVB3	114189	270084	11.71%	1.463%
52	16.368	2423	2428	2435	rVB4	67365	122717	5.32%	0.665%
53	17.686	2649	2652	2660	rVB2	56777	118864	5.16%	0.644%

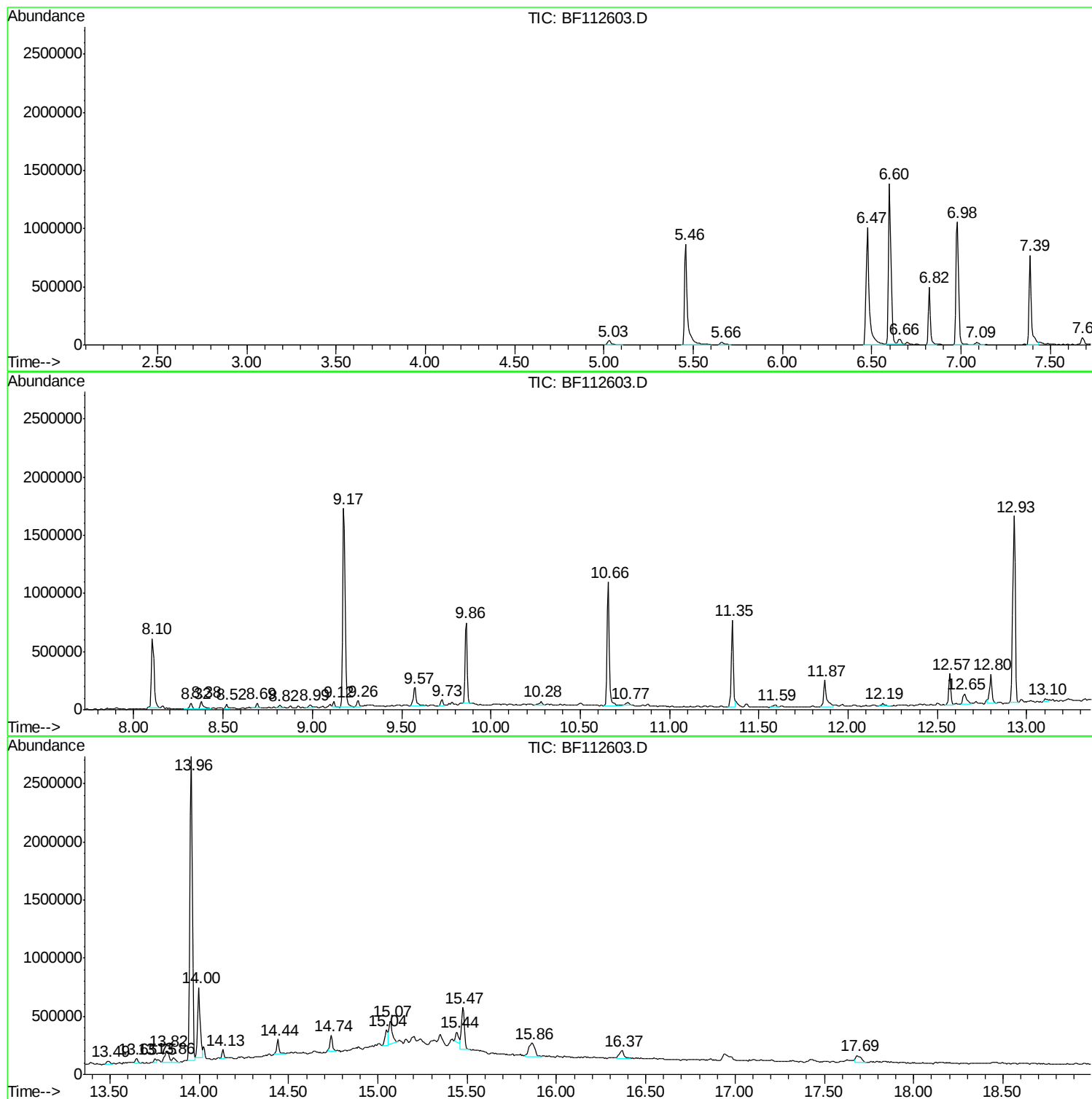
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF012519.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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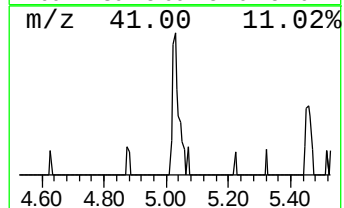
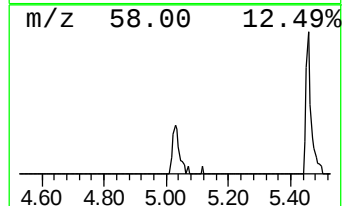
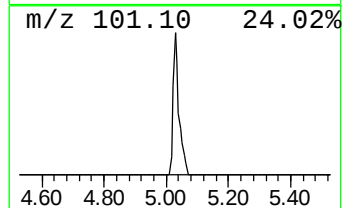
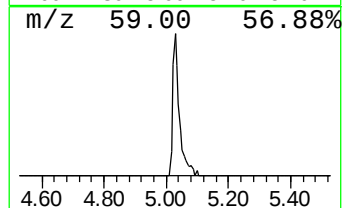
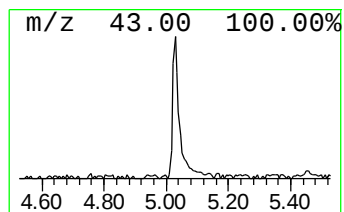
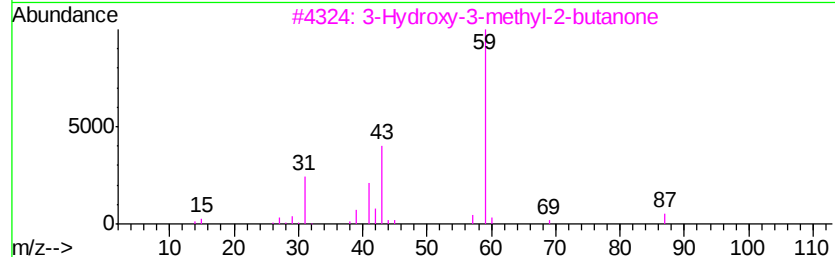
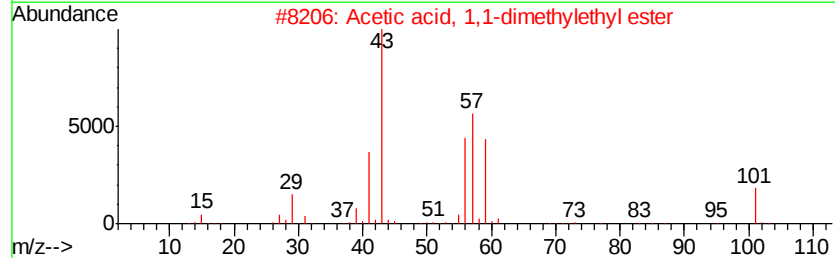
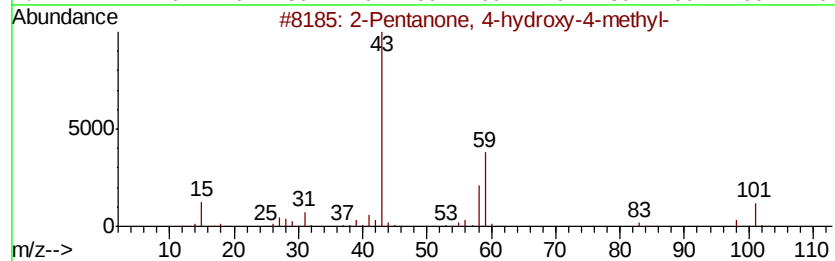
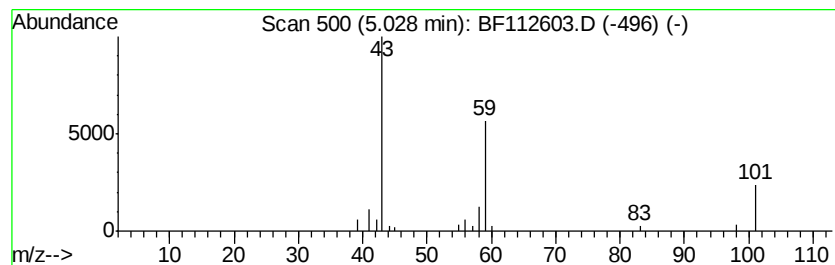
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF012519.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.03	2.74 ng	57836	1,4-Dichlorobenzene-d4	6.82

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	39
3			3-Hydroxy-3-methyl-2-butanone	102	C5H10O2	000115-22-0	23
4			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
5			5-Hexen-2-one	98	C6H10O	000109-49-9	9



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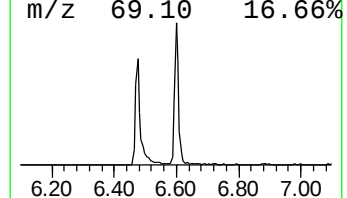
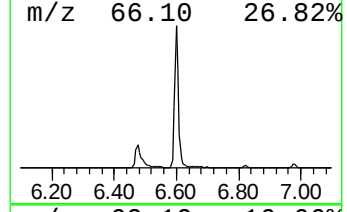
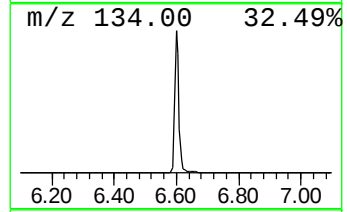
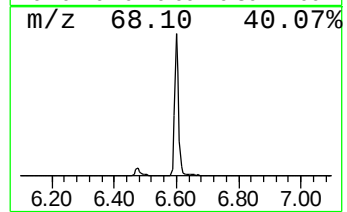
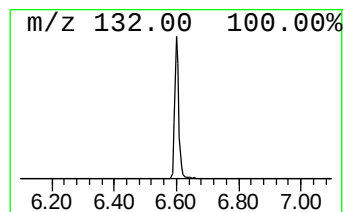
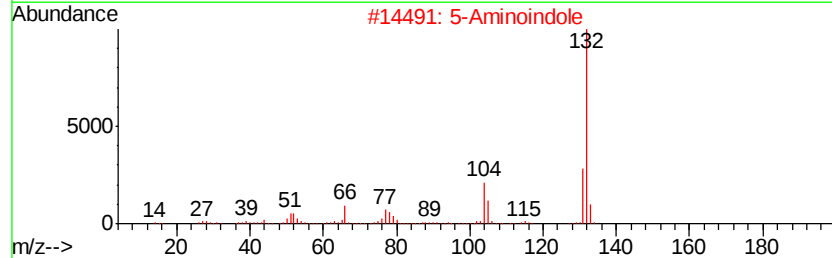
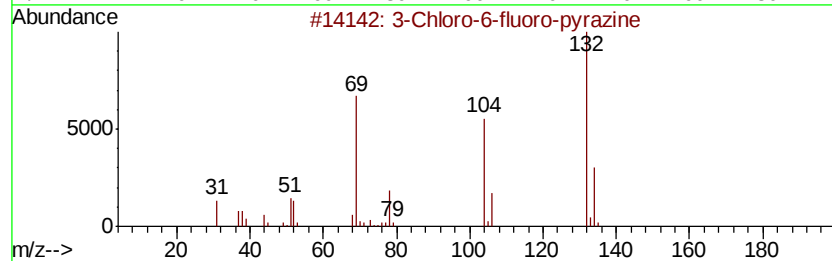
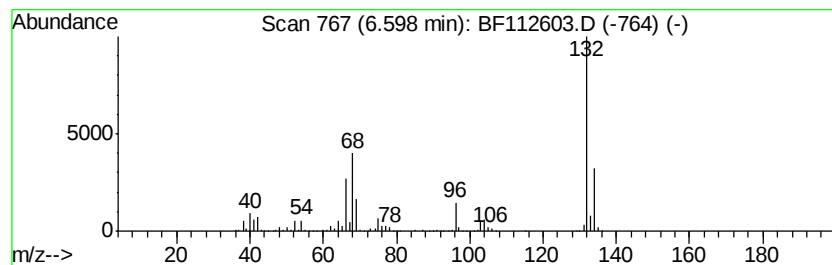
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 unknown6.60 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.60	60.65 ng	1278570	1,4-Dichlorobenzene-d4	6.82

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2			5-Aminoindole	132	C8H8N2	005192-03-0	12
3			Tranvlcyproline-propionyl	189	C12H15NO	1000123-86-3	10
4			Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
5			Benzene, 1-ethenyl-3,5-dimethyl-	132	C10H12	005379-20-4	9



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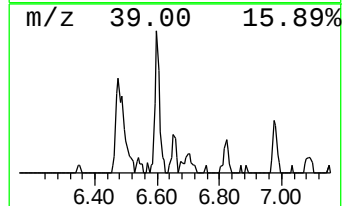
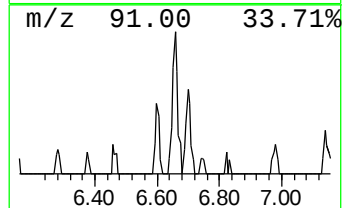
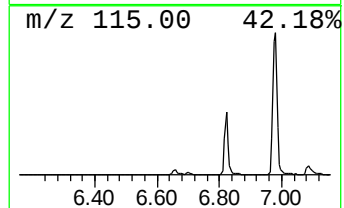
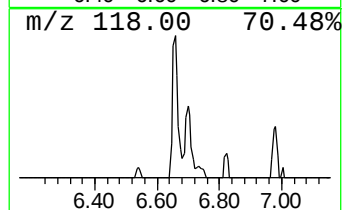
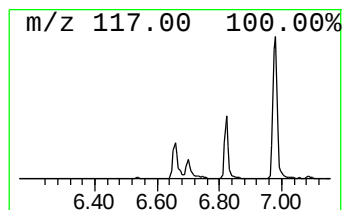
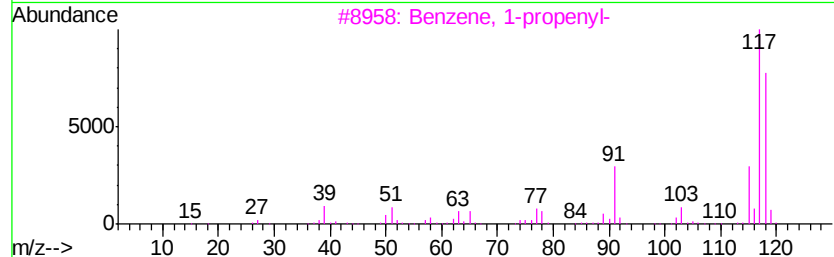
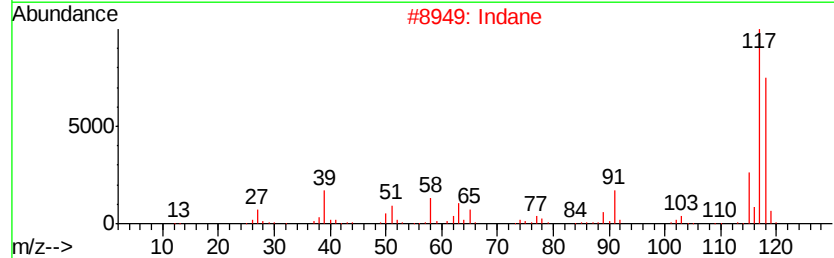
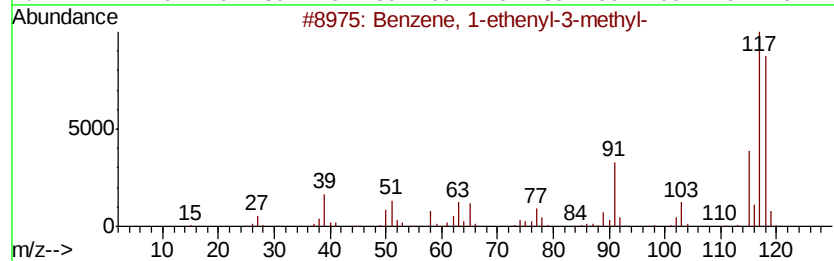
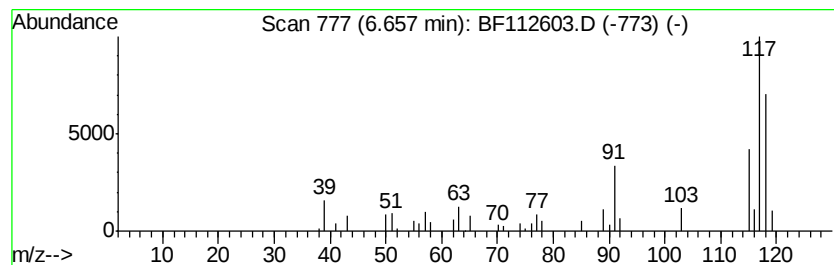
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Benzene, 1-ethenyl-3-methyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.66	2.45 ng	51745	1,4-Dichlorobenzene-d4	6.82

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethenyl-3-methyl-	118	C9H10	000100-80-1	93
2		Indane	118	C9H10	000496-11-7	81
3		Benzene, 1-propenyl-	118	C9H10	000637-50-3	81
4		Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	81
5		Benzene, cyclopropyl-	118	C9H10	000873-49-4	76



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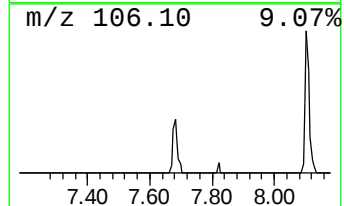
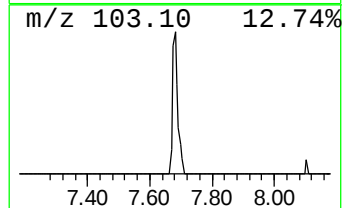
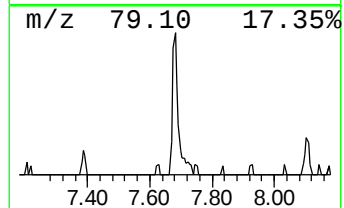
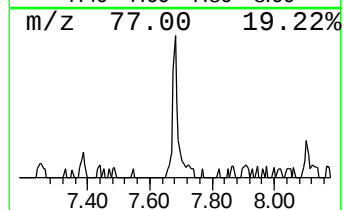
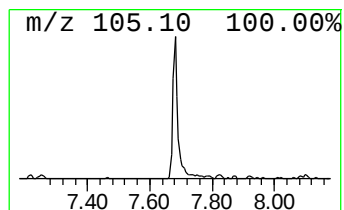
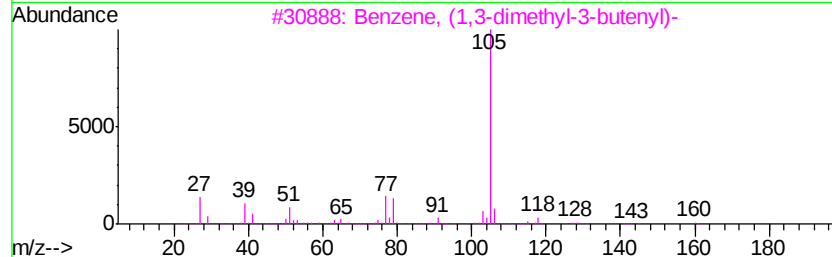
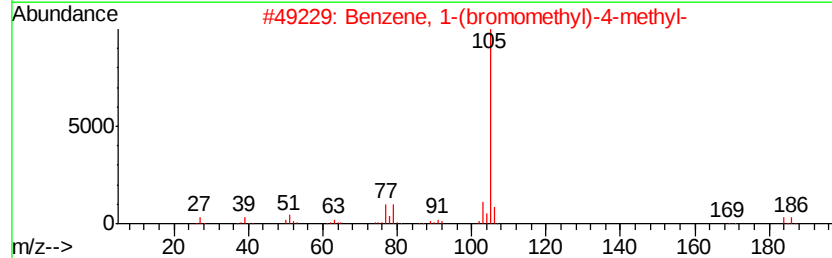
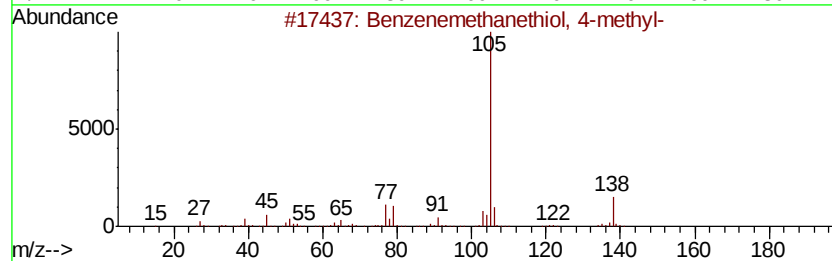
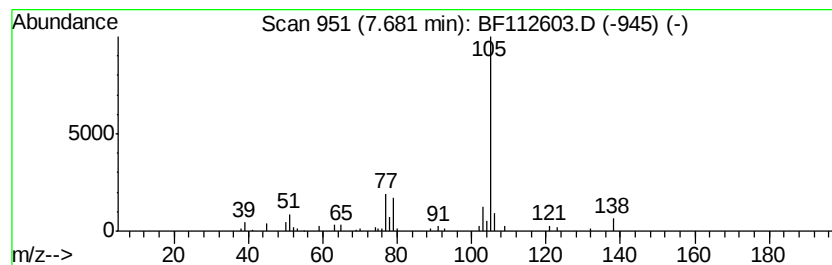
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 5 Benzenemethanethiol, 4-methyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.68	3.06 ng	80271	Naphthalene-d8	8.10

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzenemethanethiol, 4-methyl-	138	C8H10S	004498-99-1	78
2		Benzene, 1-(bromomethyl)-4-methyl-	184	C8H9Br	000104-81-4	72
3		Benzene, (1,3-dimethyl-3-butenyl)-	160	C12H16	056851-51-5	72
4		Thiophene, 2-(4-methylbenzylsulf...	252	C12H12O2S2	153837-88-8	64
5		Benzene, (1-bromoethyl)-	184	C8H9Br	000585-71-7	64



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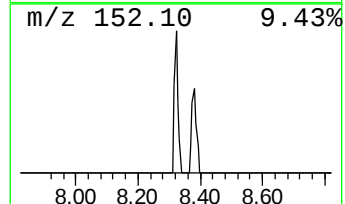
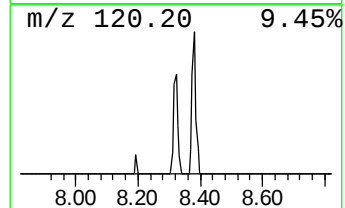
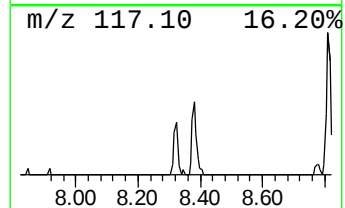
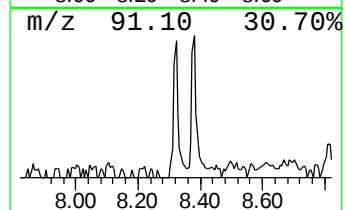
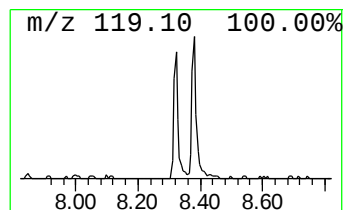
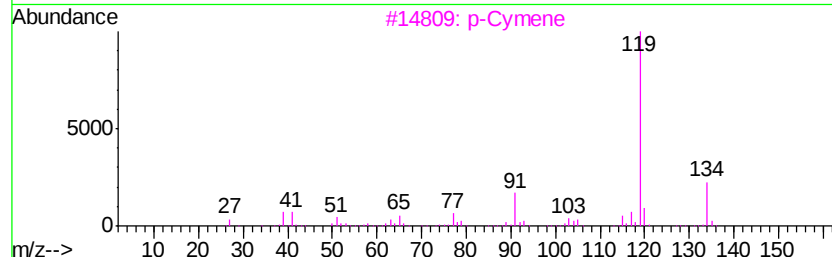
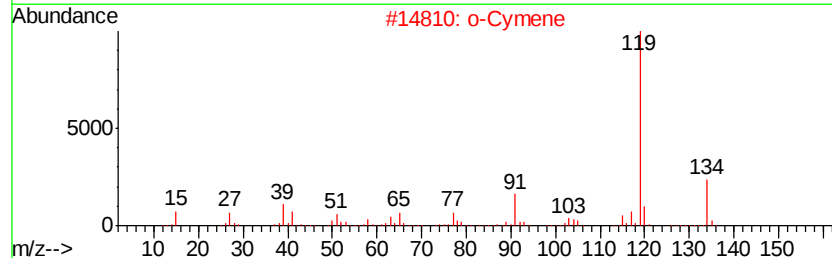
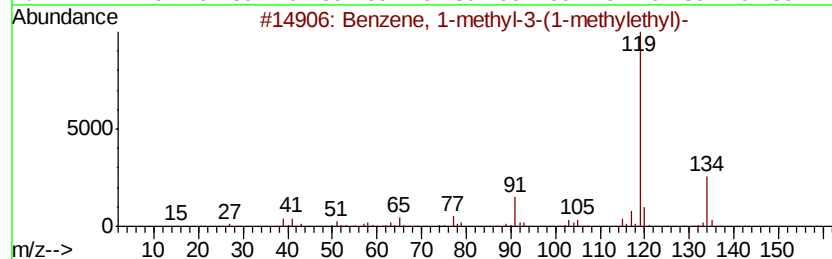
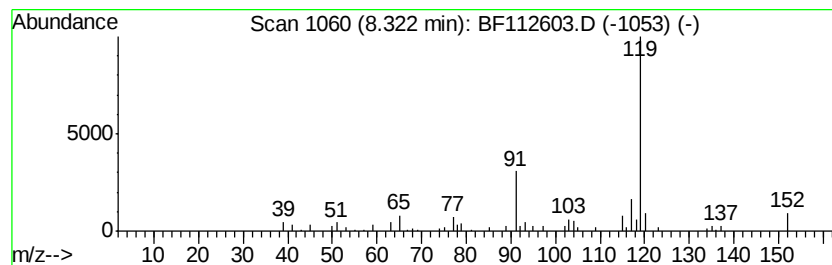
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Benzene, 1-methyl-3-(1-meth... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.32	2.51 ng	66013	Napthalene-d8	8.10

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	81
2		o-Cymene	134	C10H14	000527-84-4	81
3		p-Cymene	134	C10H14	000099-87-6	74
4		Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	72
5		Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	72



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF021819\
Data File : BF112603.D
Acq On : 18 Feb 2019 16:13
Operator : JU/SJ
Sample : K1518-15
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
CABLE-BRIDGE-FOUNDATION-D-8

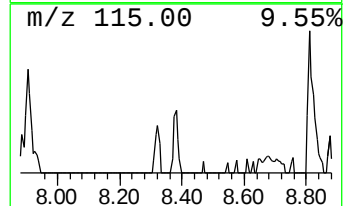
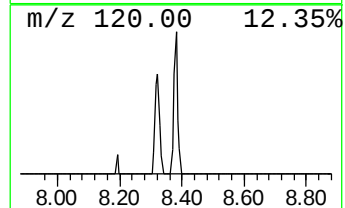
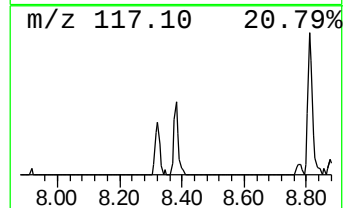
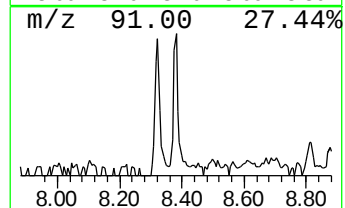
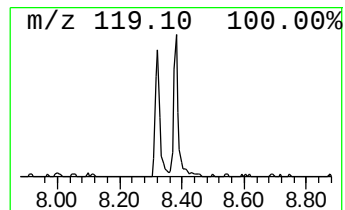
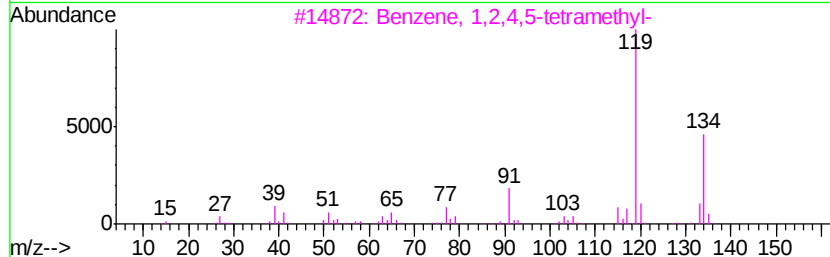
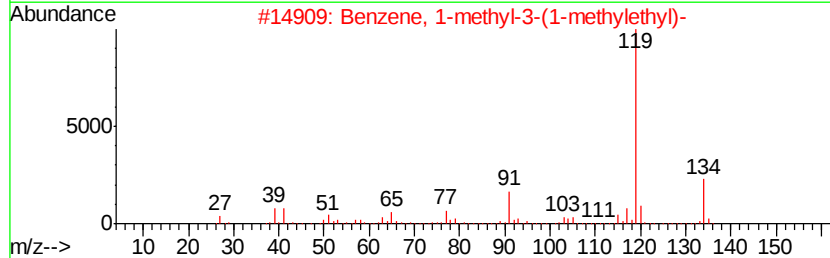
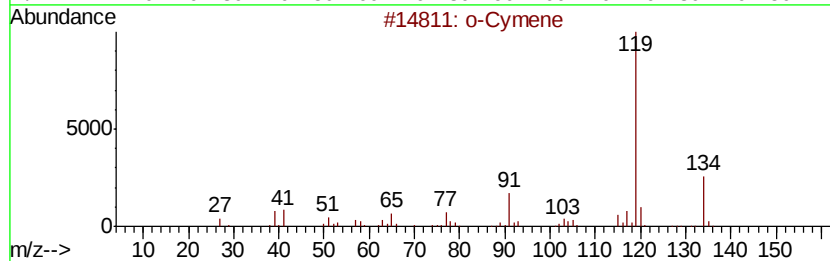
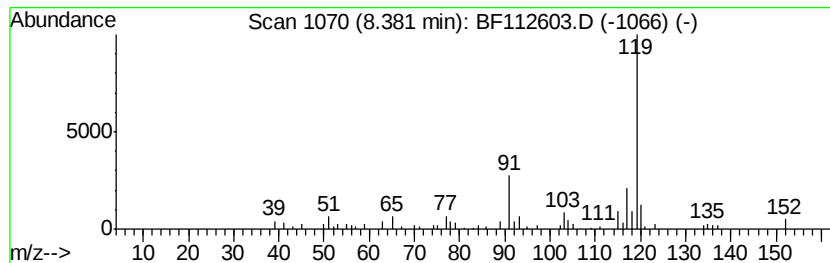
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 o-Cymene Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.38	3.13 ng	82088	Napthalene-d8	8.10

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	o-Cymene	134	C10H14	000527-84-4	81
2		Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	74
3		Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	74
4		Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	72
5		p-Cymene	134	C10H14	000099-87-6	72



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF021819\
Data File : BF112603.D
Acq On : 18 Feb 2019 16:13
Operator : JU/SJ
Sample : K1518-15
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
CABLE-BRIDGE-FOUNDATION-D-8

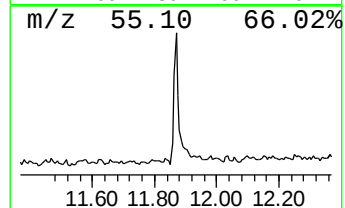
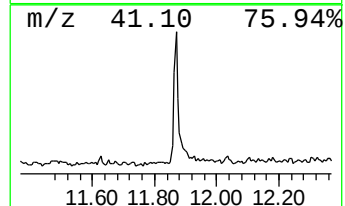
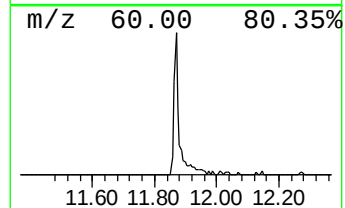
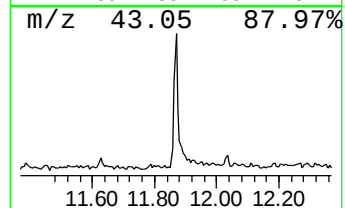
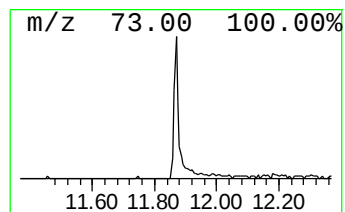
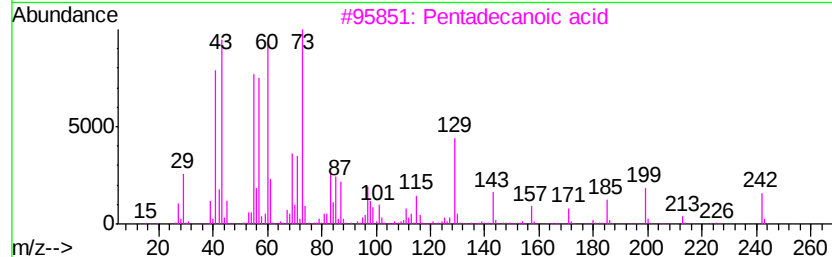
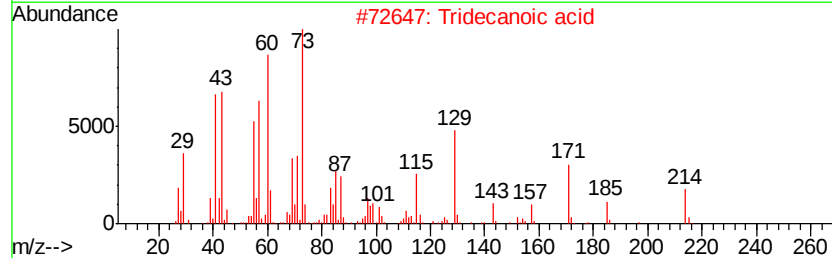
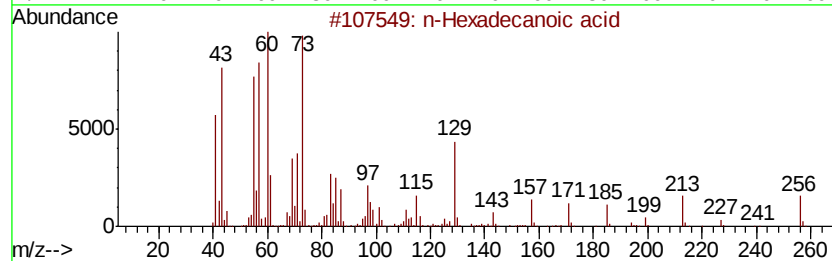
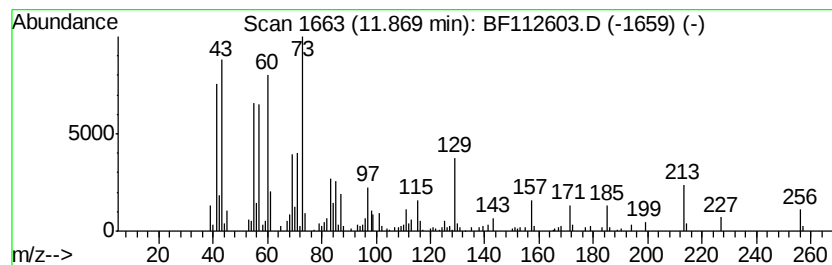
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 n-Hexadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.87	8.47 ng	277156	Phenanthrene-d10	11.35

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2		Tridecanoic acid	214	C13H26O2	000638-53-9	94
3		Pentadecanoic acid	242	C15H30O2	001002-84-2	91
4		n-Decanoic acid	172	C10H20O2	000334-48-5	76
5		1-Decanol, 2-octyl-	270	C18H38O	045235-48-1	20



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF021819\
Data File : BF112603.D
Acq On : 18 Feb 2019 16:13
Operator : JU/SJ
Sample : K1518-15
Misc :
ALS Vial : 9 Sample Multiplier: 1

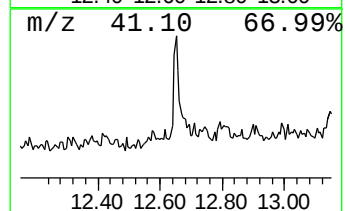
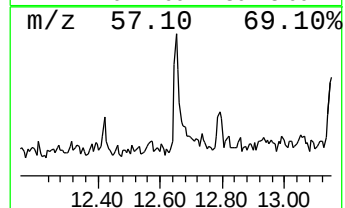
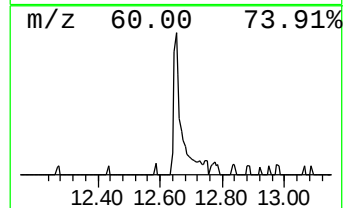
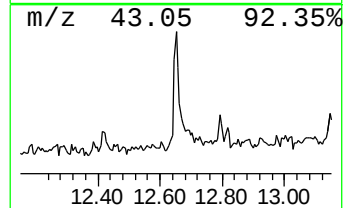
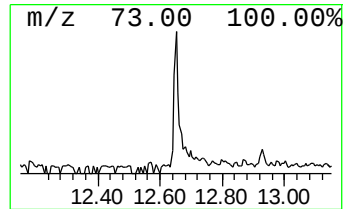
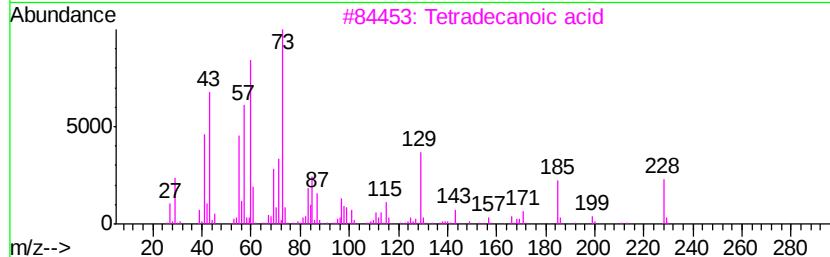
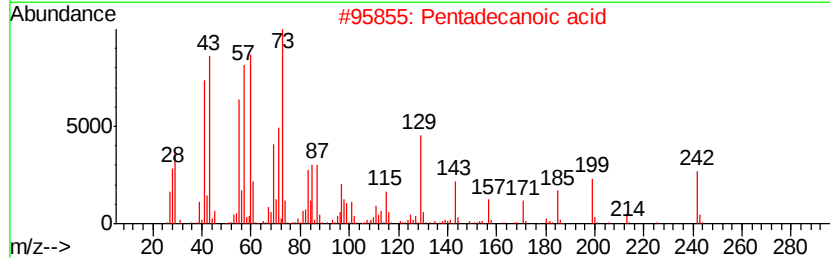
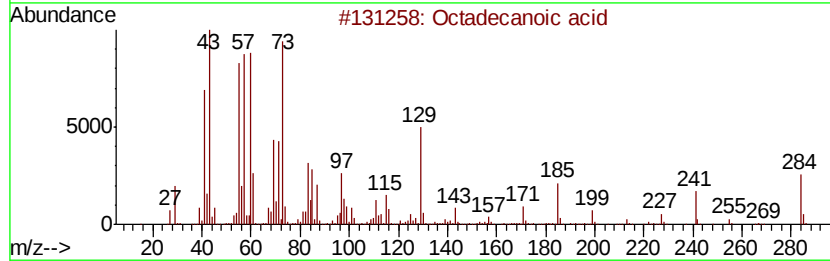
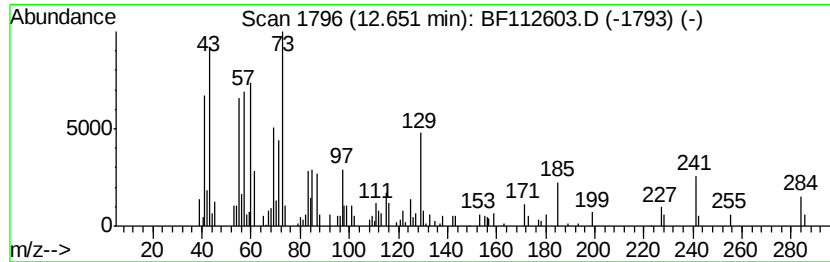
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ClientSampleId :
CABLE-BRIDGE-FOUNDATION-D-8

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF012519.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Octadecanoic acid Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD		R.T.		
12.65	3.54 ng	115906	Phenanthrene-d10		11.35		
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	90
3			Tetradecanoic acid	228	C14H28O2	000544-63-8	81
4			Tridecanoic acid	214	C13H26O2	000638-53-9	64
5			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	52



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF021819\
Data File : BF112603.D
Acq On : 18 Feb 2019 16:13
Operator : JU/SJ
Sample : K1518-15
Misc :
ALS Vial : 9 Sample Multiplier: 1

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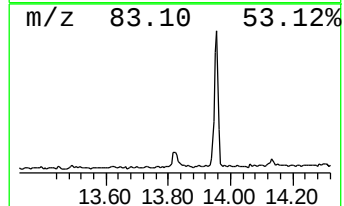
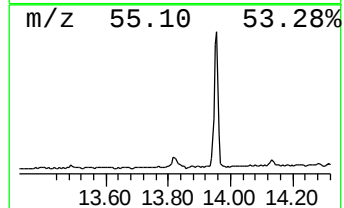
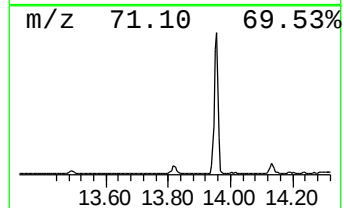
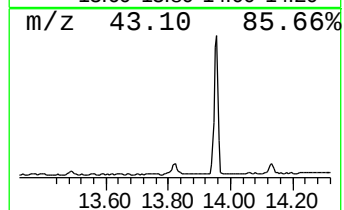
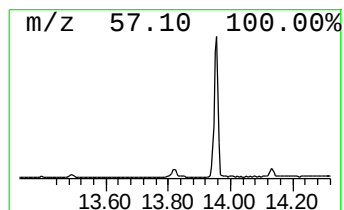
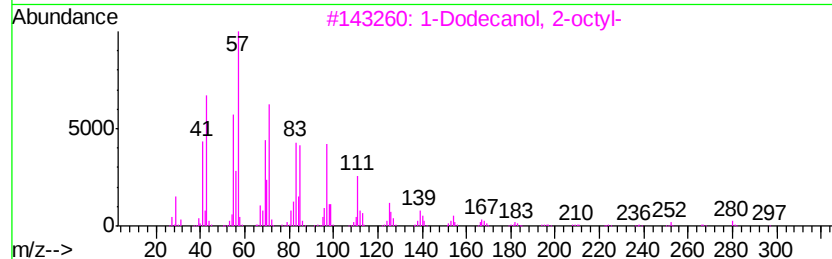
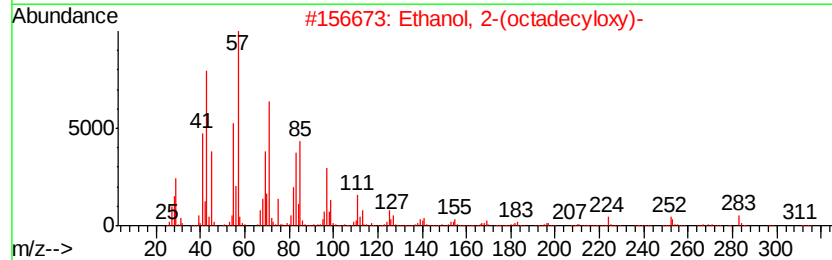
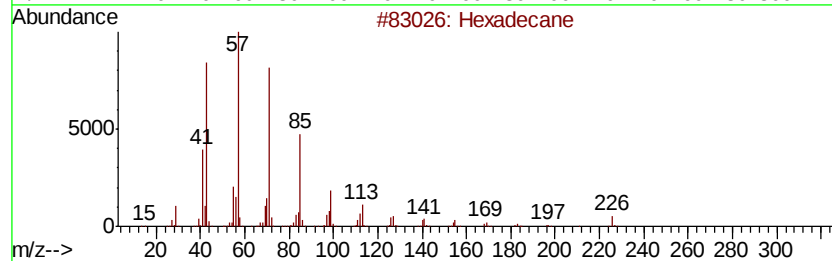
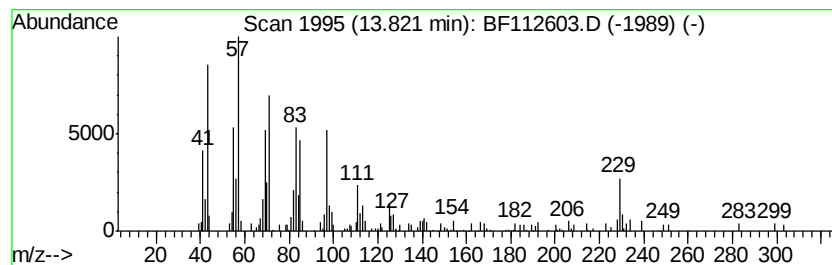
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Hexadecane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.82	5.44 ng	164314	Chrysene-d12	14.00

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexadecane	226	C16H34	000544-76-3	86
2			Ethanol, 2-(octadecyloxy)-	314	C20H42O2	002136-72-3	81
3			1-Dodecanol, 2-octyl-	298	C20H42O	005333-42-6	68
4			Sulfurous acid, octadecyl 2-prop...	376	C21H44O3S	1000309-12-7	62
5			Cetene	224	C16H32	000629-73-2	55



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF021819\
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Sample : K1518-15
Misc :
ALS Vial : 9 Sample Multiplier: 1

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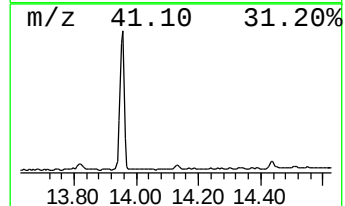
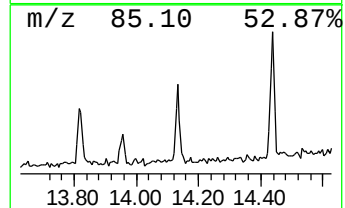
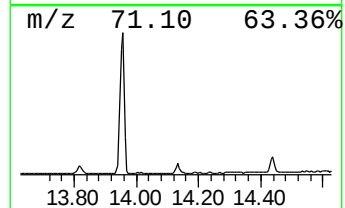
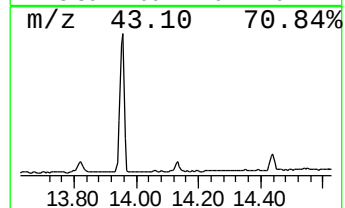
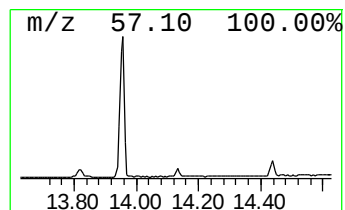
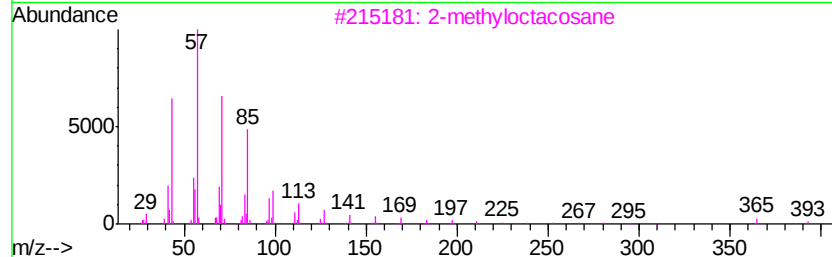
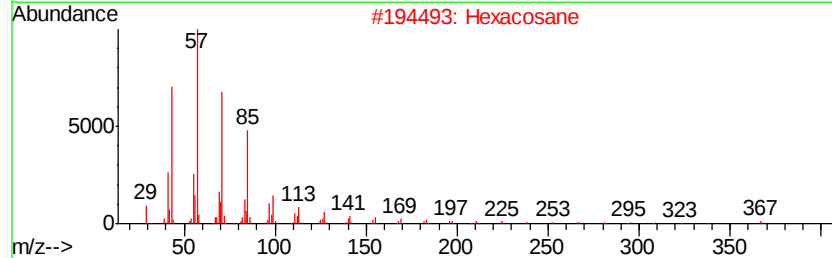
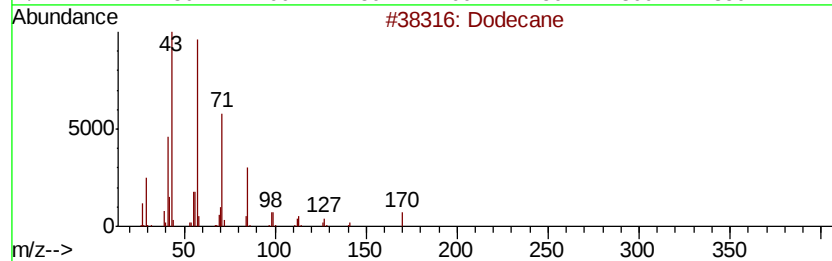
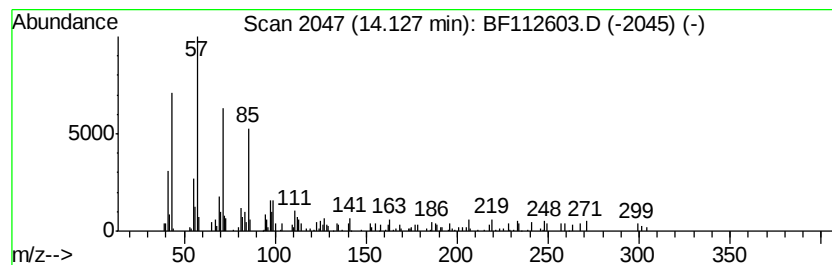
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Dodecane Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.13	2.12 ng	64197	Chrysene-d12	14.00

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Dodecane	170	C12H26	000112-40-3	86
2		Hexacosane	366	C26H54	000630-01-3	76
3		2-methyloctacosane	408	C29H60	1000376-72-8	68
4		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	68
5		Tetratetracontane	619	C44H90	007098-22-8	68



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF021819\
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ALS Vial : 9 Sample Multiplier: 1

Instrument :
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ClientSampleId :
CABLE-BRIDGE-FOUNDATION-D-8

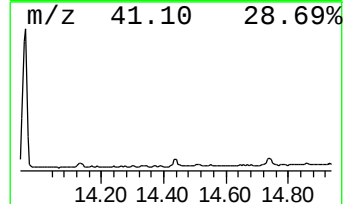
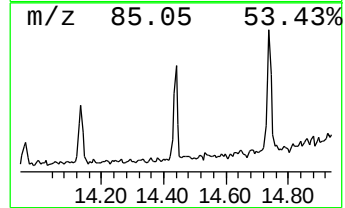
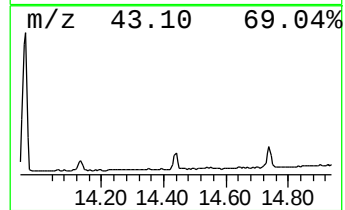
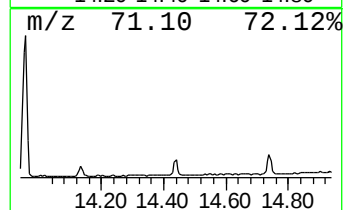
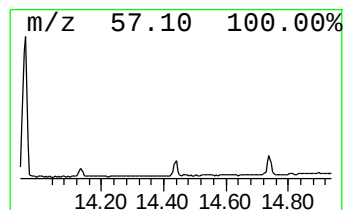
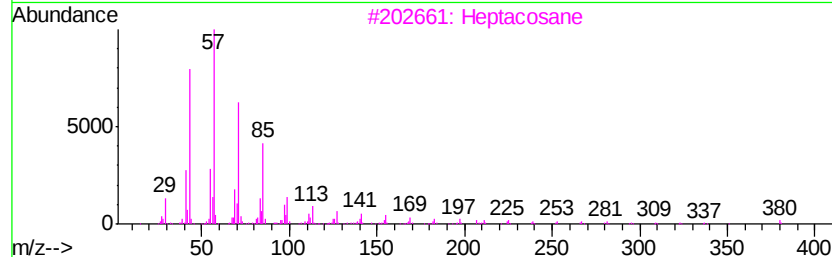
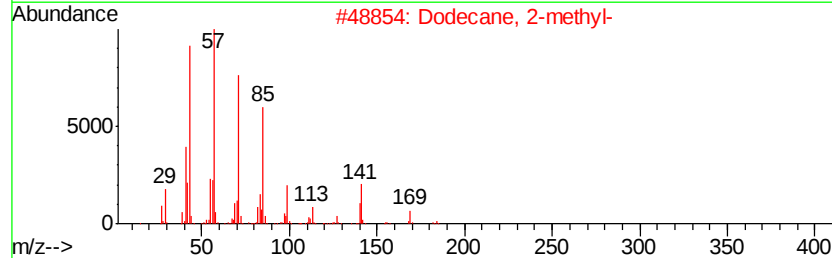
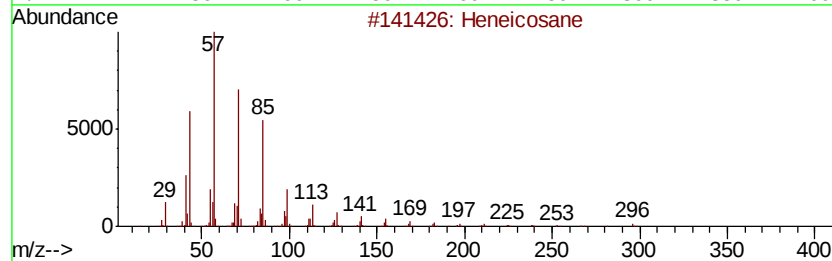
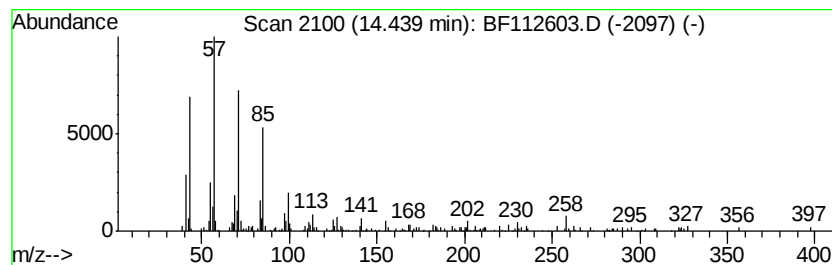
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Heneicosane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.44	4.21 ng	127143	Chrysene-d12	14.00

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosane	296	C21H44	000629-94-7	87
2		Dodecane, 2-methyl-	184	C13H28	001560-97-0	83
3		Heptacosane	380	C27H56	000593-49-7	81
4		Hexacosane	366	C26H54	000630-01-3	81
5		Tetracosane	338	C24H50	000646-31-1	81



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF021819\
Data File : BF112603.D
Acq On : 18 Feb 2019 16:13
Operator : JU/SJ
Sample : K1518-15
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
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ClientSampleId :
CABLE-BRIDGE-FOUNDATION-D-8

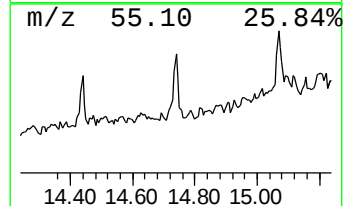
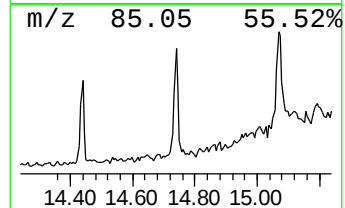
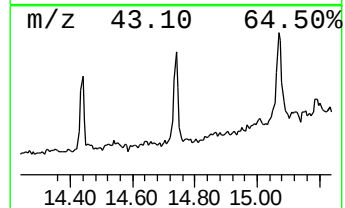
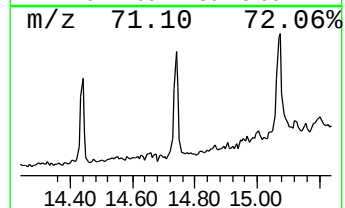
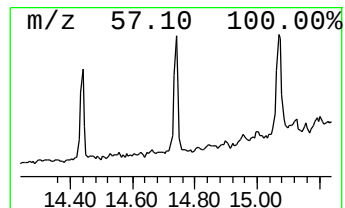
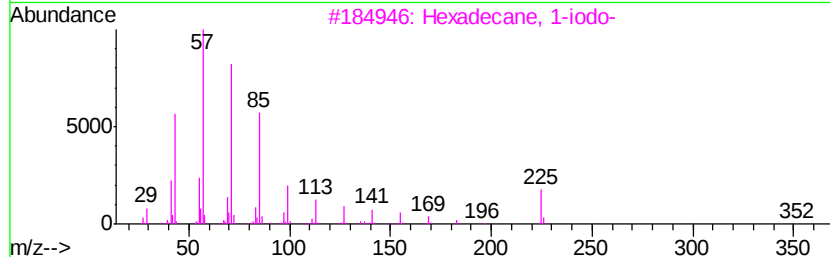
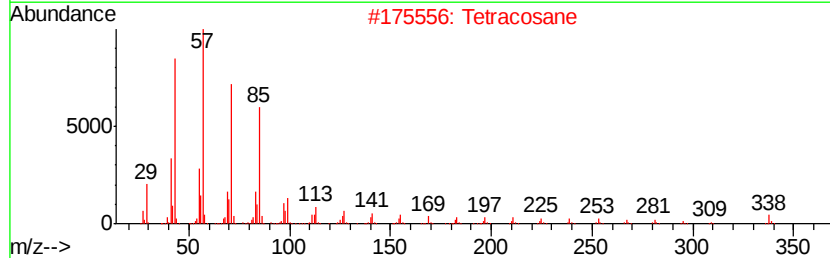
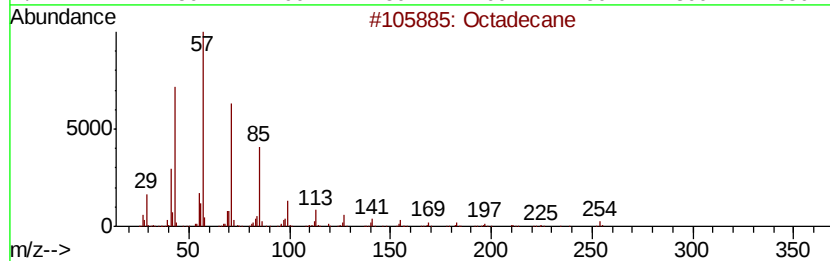
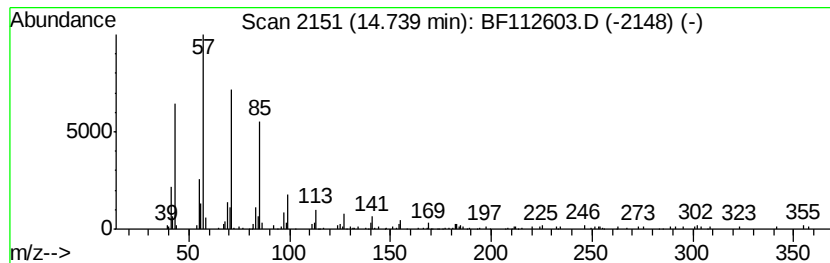
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Octadecane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.74	5.54 ng	128647	Perylene-d12	15.47

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecane	254	C18H38	000593-45-3	90
2		Tetracosane	338	C24H50	000646-31-1	87
3		Hexadecane, 1-iodo-	352	C16H33I	000544-77-4	87
4		Hexadecane	226	C16H34	000544-76-3	87
5		Hexacosane	366	C26H54	000630-01-3	86



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF021819\
Data File : BF112603.D
Acq On : 18 Feb 2019 16:13
Operator : JU/SJ
Sample : K1518-15
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
CABLE-BRIDGE-FOUNDATION-D-8

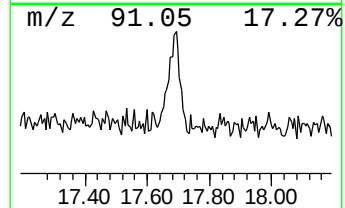
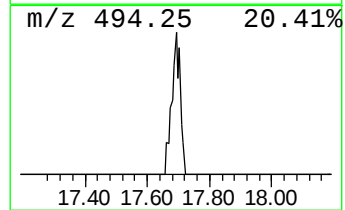
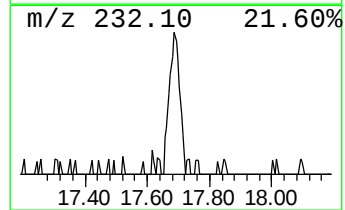
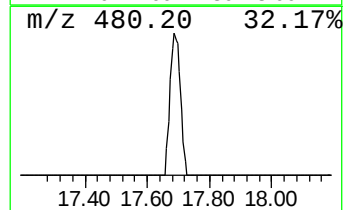
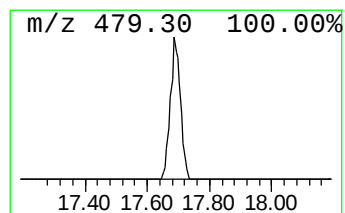
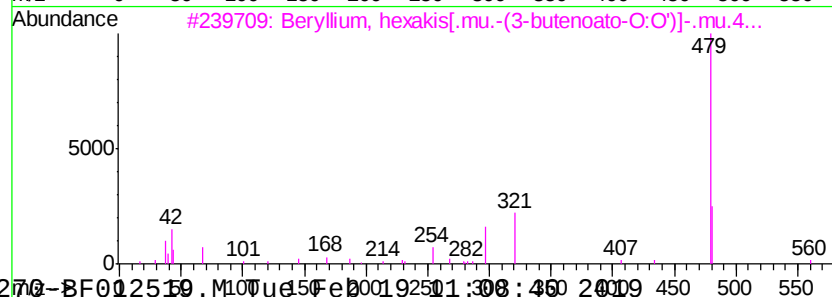
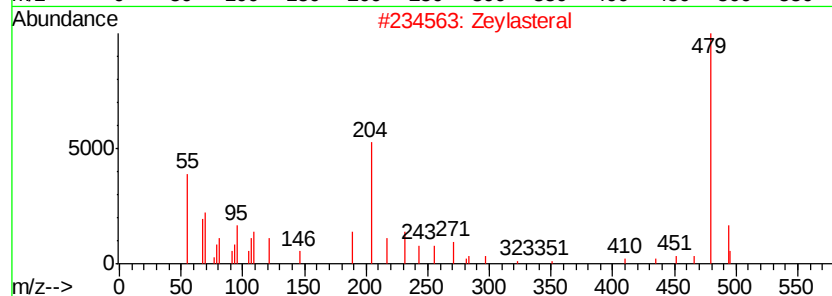
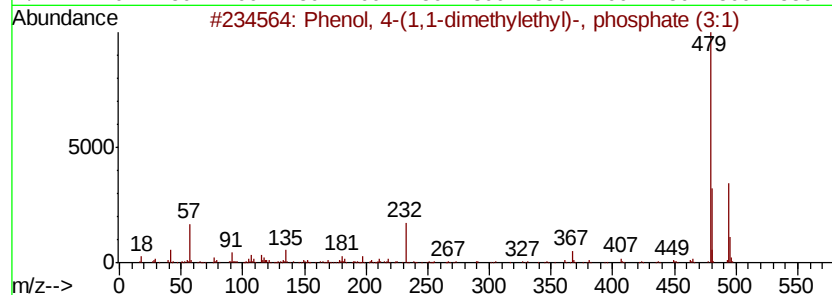
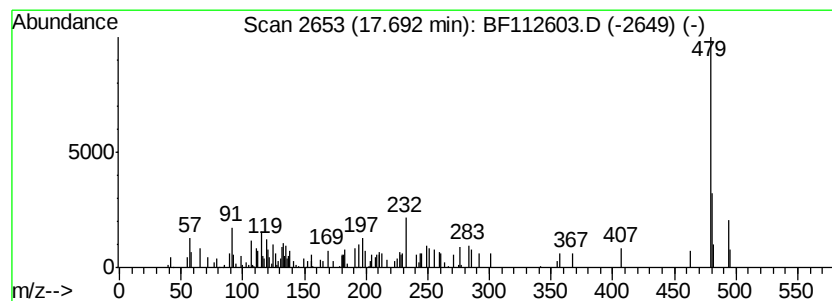
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 unknown17.69 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.69	5.12 ng	118864	Perylene-d12	15.47

Hit#	of	4	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenol, 4-(1,1-dimethylethyl)-, ...	494	C30H39O4P	000078-33-1	38
2			Zevlasteral	494	C30H38O6	087064-16-2	38
3			Beryllium, hexakis[.mu.-(3-buten...	562	C24H30Be4O13	056377-88-9	9
4			Silane, diethyl(dodec-9-ynyloxy)...	508	C32H64O2Si	1000363-58-8	2



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF021819\
 Data File : BF112603.D
 Acq On : 18 Feb 2019 16:13
 Operator : JU/SJ
 Sample : K1518-15
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CABLE-BRIDGE-FOUNDATION-D-8

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF012519.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
2-Pentanone, 4-hy...	5.03	2.7	ng	57836	1	6.82	421603	20.0
unknown6.60	6.60	60.6	ng	1278570	1	6.82	421603	20.0
Benzene, 1-etheny...	6.66	2.5	ng	51745	1	6.82	421603	20.0
Benzenemethanethi...	7.68	3.1	ng	80271	2	8.10	525119	20.0
Benzene, 1-methyl...	8.32	2.5	ng	66013	2	8.10	525119	20.0
o-Cymene	8.38	3.1	ng	82088	2	8.10	525119	20.0
n-Hexadecanoic acid	11.87	8.5	ng	277156	4	11.35	654606	20.0
Octadecanoic acid	12.65	3.5	ng	115906	4	11.35	654606	20.0
Hexadecane	13.82	5.4	ng	164314	5	14.00	604307	20.0
Dodecane	14.13	2.1	ng	64197	5	14.00	604307	20.0
Heneicosane	14.44	4.2	ng	127143	5	14.00	604307	20.0
Octadecane	14.74	5.5	ng	128647	6	15.47	464501	20.0
unknown17.69	17.69	5.1	ng	118864	6	15.47	464501	20.0