

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF041020\
 Data File : BF119910.D
 Acq On : 10 Apr 2020 22:15
 Operator : MA/SJ
 Sample : L2199-05
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-39-38-COMP

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-BF040820.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	3.751	278	283	289	rBV2	26676	37068	1.14%	0.120%
2	4.946	482	486	497	rVB	401597	465303	14.37%	1.510%
3	5.363	553	557	561	rBV	3302170	3238568	100.00%	10.512%
4	6.387	726	731	743	rVB	2848041	3134010	96.77%	10.173%
5	6.581	761	764	770	rBV	197071	177173	5.47%	0.575%
6	6.751	790	793	797	rBV	1161770	1034915	31.96%	3.359%
7	6.910	816	820	823	rBV	3014980	2483407	76.68%	8.061%
8	7.316	884	889	892	rBV	2149685	1946484	60.10%	6.318%
9	8.039	1007	1012	1015	rBV	1553870	1262512	38.98%	4.098%
10	8.634	1110	1113	1117	rBV	39763	40186	1.24%	0.130%
11	9.116	1190	1195	1198	rBV	3593501	3116949	96.24%	10.117%
12	9.204	1206	1210	1212	rBV	56243	66342	2.05%	0.215%
13	9.239	1212	1216	1219	rVV	1416859	1211307	37.40%	3.932%
14	9.510	1259	1262	1266	rVB	475996	407117	12.57%	1.321%
15	9.604	1275	1278	1285	rBV	250820	311135	9.61%	1.010%
16	9.657	1285	1287	1293	rVB	46653	50684	1.57%	0.165%
17	9.733	1297	1300	1304	rVB	46767	41765	1.29%	0.136%
18	9.792	1304	1310	1314	rBV	1311787	1200101	37.06%	3.895%
19	10.175	1372	1375	1377	rBV	44349	36319	1.12%	0.118%
20	10.380	1407	1410	1417	rBV	291140	308288	9.52%	1.001%
21	10.580	1440	1444	1448	rBV	1803116	1647719	50.88%	5.348%
22	10.622	1448	1451	1457	rVB5	26608	42540	1.31%	0.138%
23	10.710	1462	1466	1474	rVB4	61133	93467	2.89%	0.303%
24	11.186	1544	1547	1552	rVB2	32515	41780	1.29%	0.136%
25	11.280	1559	1563	1566	rBV	1329076	1078552	33.30%	3.501%
26	11.304	1566	1567	1570	rVV	78243	44133	1.36%	0.143%
27	11.357	1571	1576	1578	rVB	45368	50724	1.57%	0.165%
28	11.475	1592	1596	1600	rBV	48155	62853	1.94%	0.204%
29	11.527	1601	1605	1611	rVB8	23348	40876	1.26%	0.133%
30	11.816	1650	1654	1660	rBV2	714442	665125	20.54%	2.159%
31	11.892	1664	1667	1670	rBV	54110	55063	1.70%	0.179%
32	12.345	1739	1744	1748	rBV2	116887	222962	6.88%	0.724%
33	12.422	1754	1757	1761	rBV5	31807	45621	1.41%	0.148%
34	12.492	1766	1769	1772	rBV	121339	123666	3.82%	0.401%

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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	12.533	1773	1776	1780	rVB	160652	161665	4.99%	0.525%
36	12.604	1784	1788	1795	rBV2	405847	473620	14.62%	1.537%
37	12.721	1805	1808	1811	rBV	209973	176815	5.46%	0.574%
38	12.751	1811	1813	1816	rVB2	65016	56619	1.75%	0.184%
39	12.869	1829	1833	1837	rBV	2579473	2448022	75.59%	7.946%
40	12.945	1843	1846	1850	rBV2	41515	62887	1.94%	0.204%
41	13.280	1901	1903	1906	rVB	52528	39450	1.22%	0.128%
42	13.786	1985	1989	1996	rVB2	226789	348193	10.75%	1.130%
43	13.921	2008	2012	2015	rBV	1096377	1065851	32.91%	3.460%
44	14.404	2091	2094	2096	rBV2	158790	138512	4.28%	0.450%
45	14.680	2139	2141	2143	rBV	170650	151164	4.67%	0.491%
46	15.033	2198	2201	2208	rVB2	179561	231651	7.15%	0.752%
47	15.362	2253	2257	2260	rBV	564537	669397	20.67%	2.173%

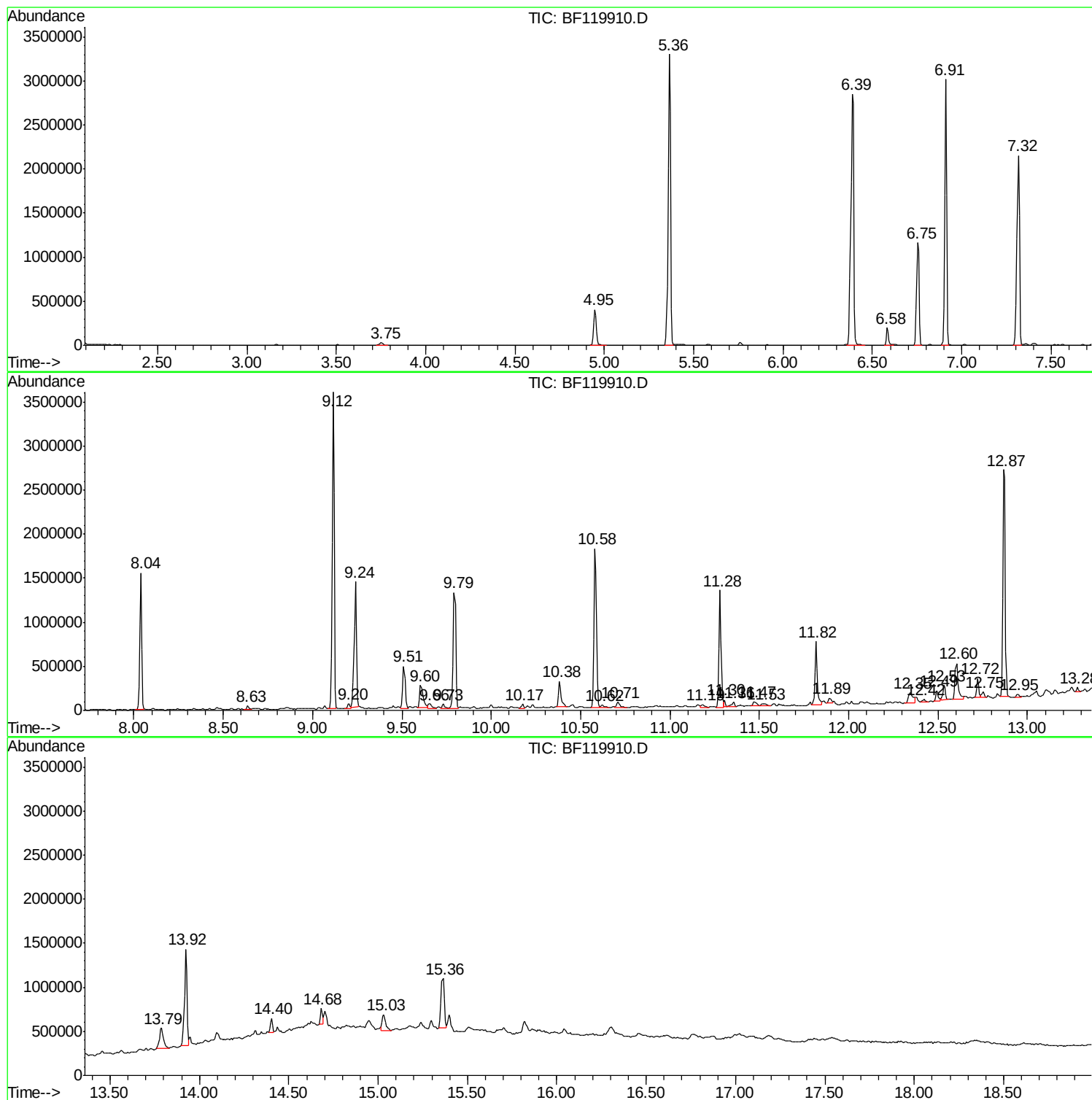
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF040820.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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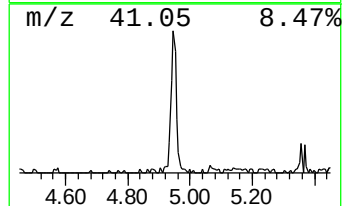
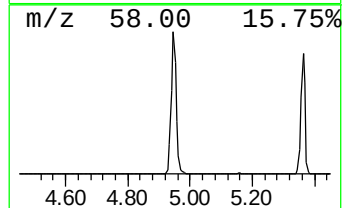
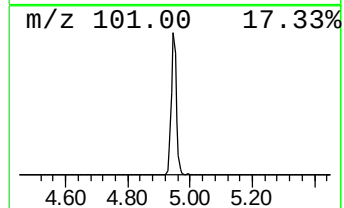
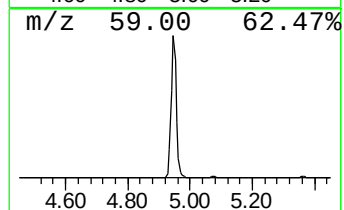
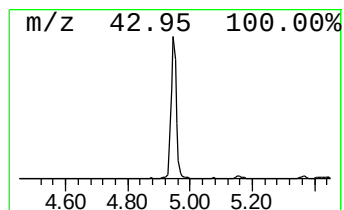
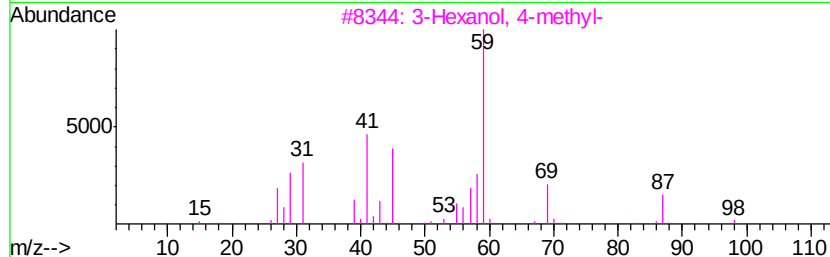
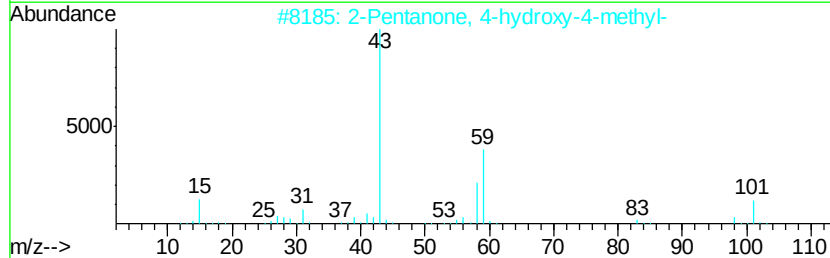
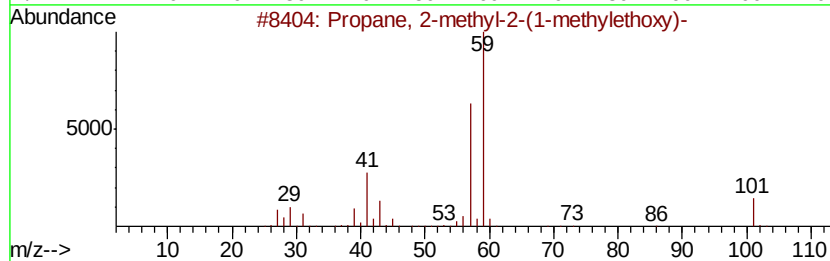
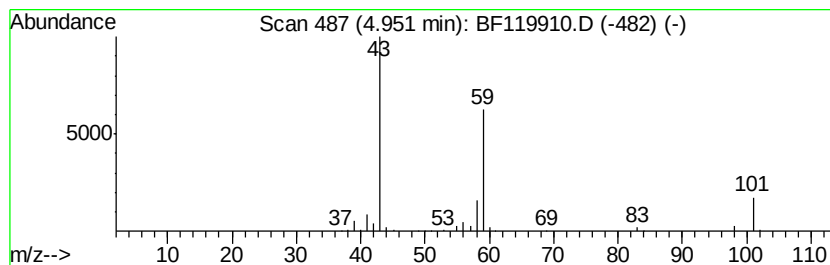
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF040820.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Propane, 2-methyl-2-(1-meth... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.95	8.99 ng	465303	1,4-Dichlorobenzene-d4	6.76

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	53
2		2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
3		3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33
4		2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	28
5		Guanidine	59	CH5N3	000113-00-8	9



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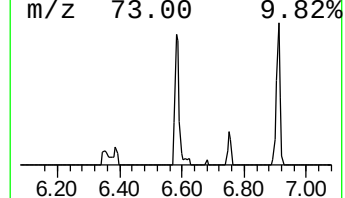
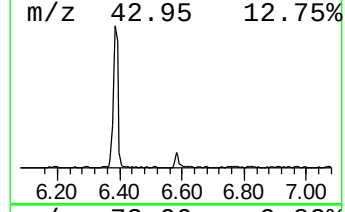
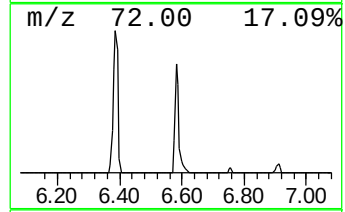
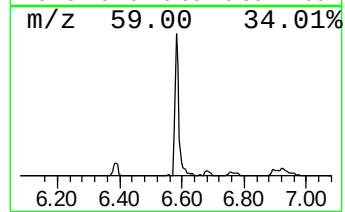
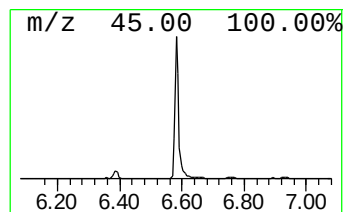
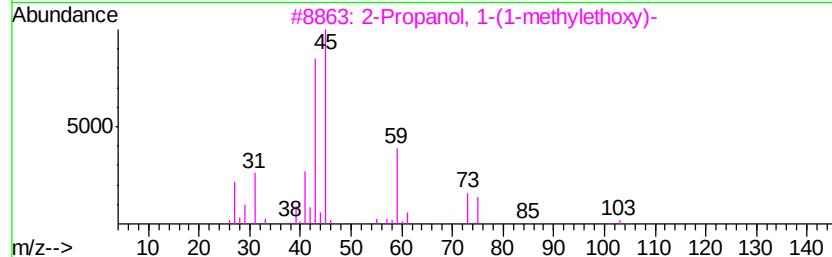
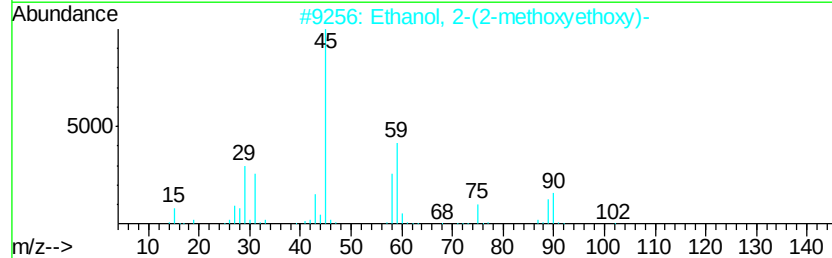
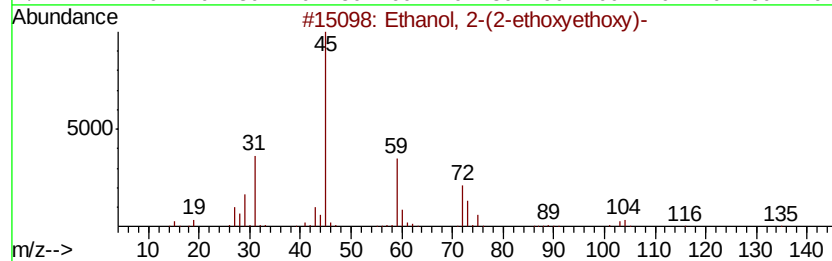
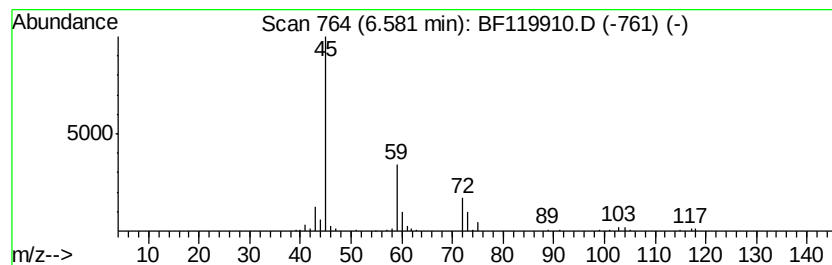
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Ethanol, 2-(2-ethoxyethoxy)- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.58	3.42 ng	177173	1,4-Dichlorobenzene-d4	6.76

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Ethanol, 2-(2-ethoxyethoxy)-	134	C6H14O3	000111-90-0	50
2		Ethanol, 2-(2-methoxyethoxy)-	120	C5H12O3	000111-77-3	45
3		2-Propanol, 1-(1-methylethoxy)-	118	C6H14O2	003944-36-3	36
4		Ethanol, 2-[2-(2-propenyloxy)eth...]	146	C7H14O3	015075-50-0	33
5		Methoxyacetic acid, butyl ester	146	C7H14O3	017640-22-1	33



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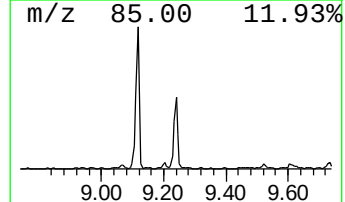
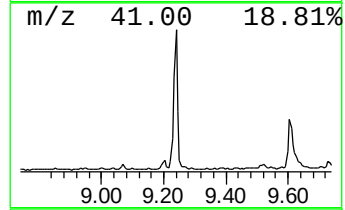
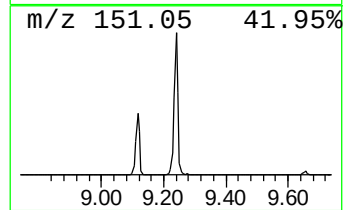
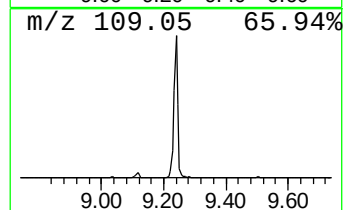
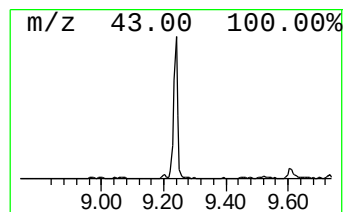
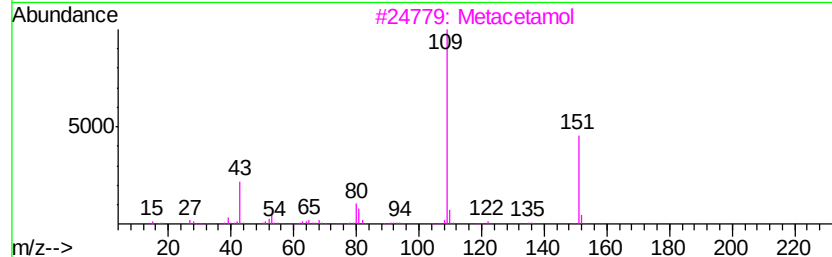
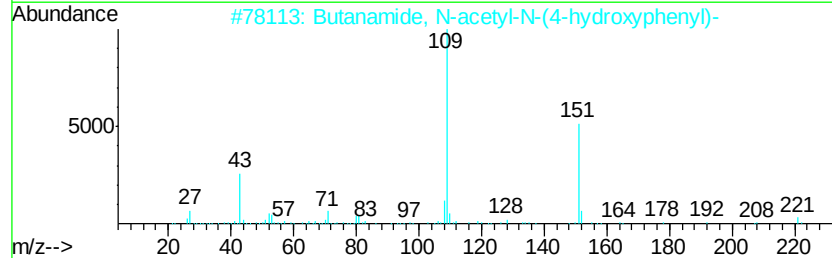
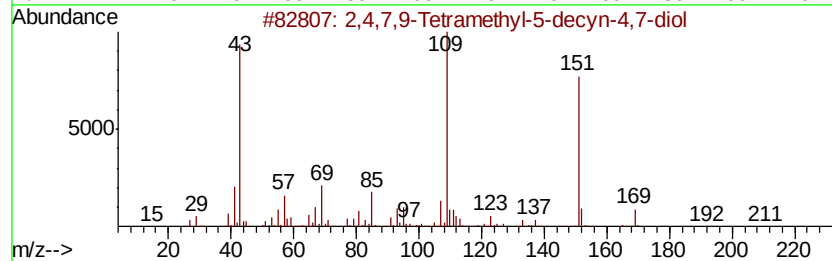
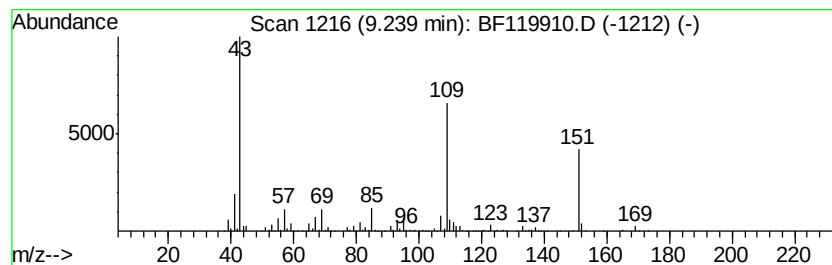
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.	
9.24	20.19 ng	1211310	Acenaphthene-d10	9.79	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,4,7,9-Tetramethyl-5-decyn-4,7-...	226	C14H26O2	000126-86-3	87
2	Butanamide, N-acetyl-N-(4-hydrox...	221	C12H15NO3	1000107-08-7	42
3	Metacetamol	151	C8H9NO2	000621-42-1	42
4	Cyclohexanol, 3,3,5-trimethyl-	142	C9H18O	000116-02-9	38
5	Pyridine, 3-butyl-, 1-oxide	151	C9H13NO	031396-33-5	36



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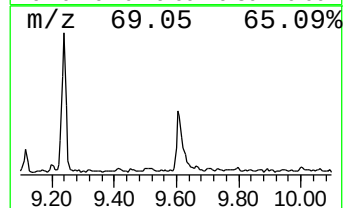
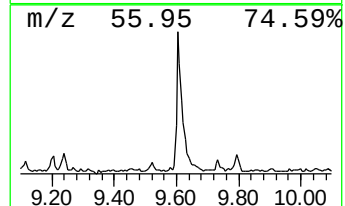
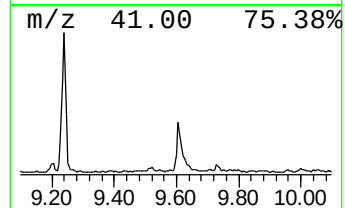
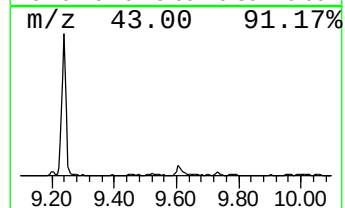
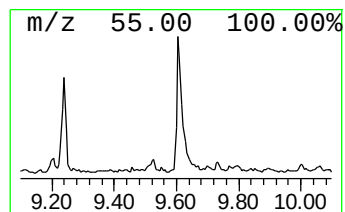
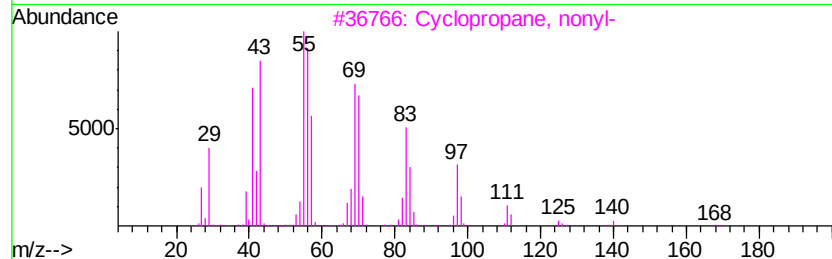
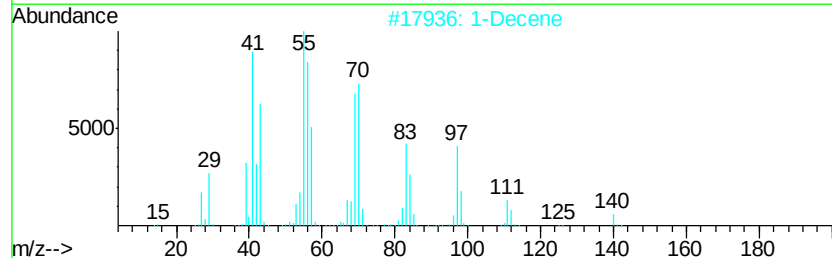
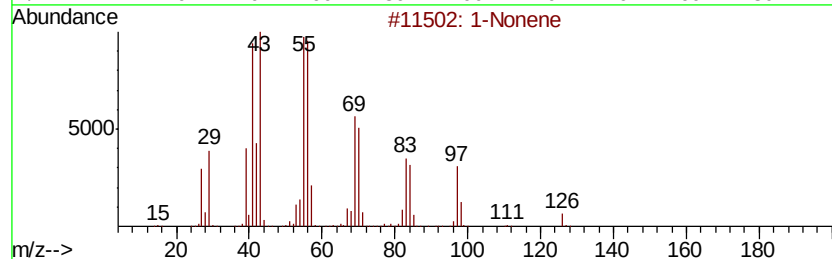
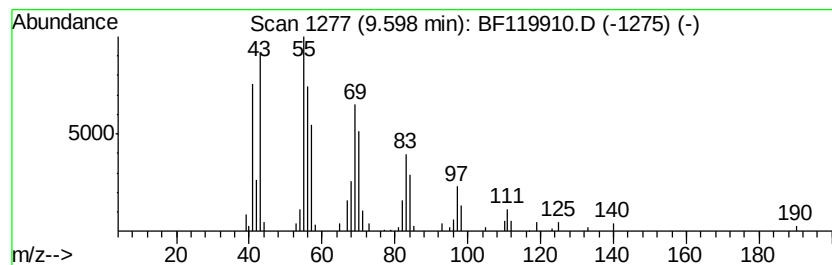
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Peak Number 5 1-Nonene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.60	5.19 ng	311135	Acenaphthene-d10	9.79

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Nonene		126	C9H18	000124-11-8	94
2	1-Decene		140	C10H20	000872-05-9	93
3	Cyclopropane, nonyl-		168	C12H24	074663-85-7	91
4	1-Undecanol		172	C11H24O	000112-42-5	91
5	Cyclooctane, 1,2-dimethyl-		140	C10H20	013151-94-5	90



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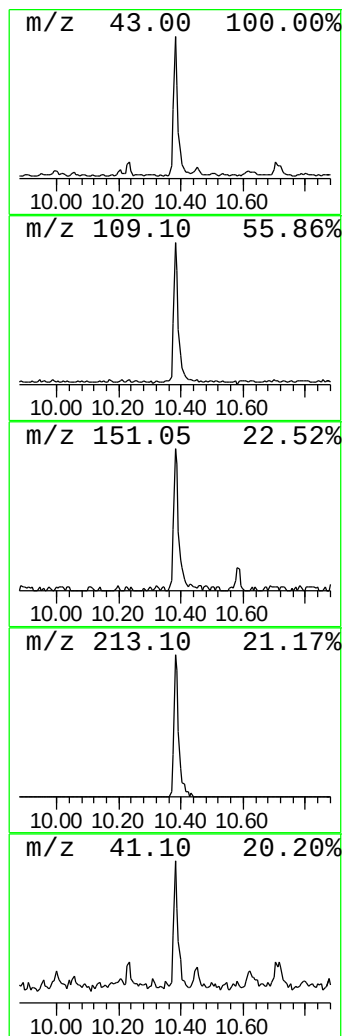
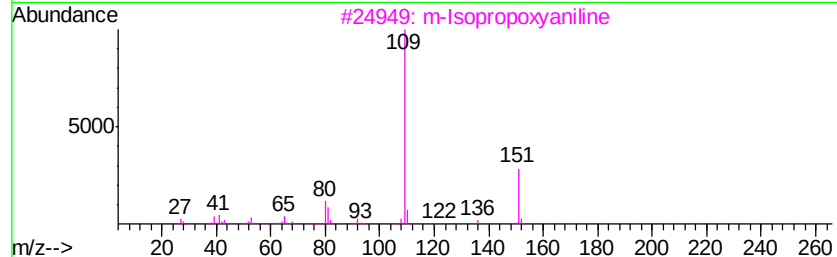
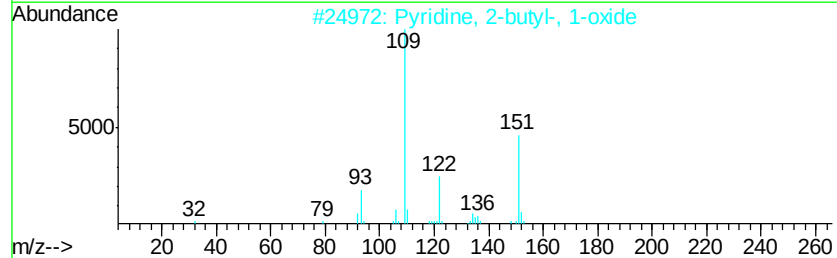
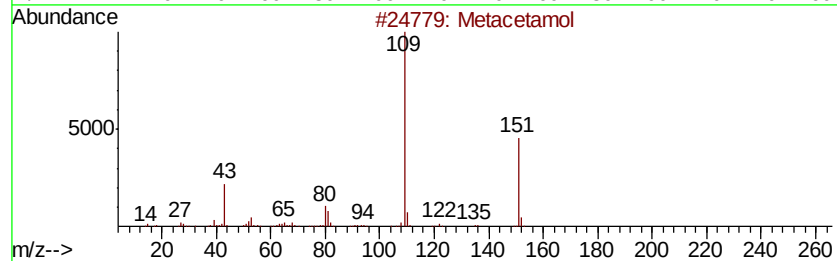
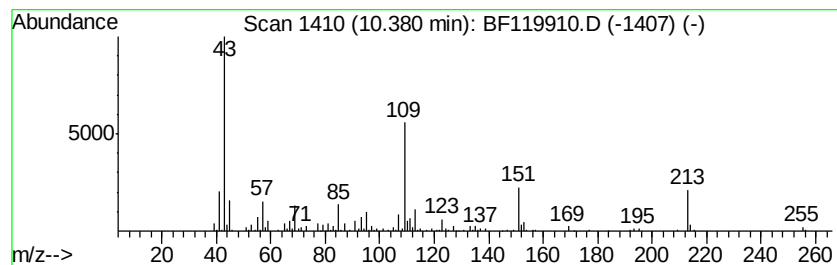
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Peak Number 6 unknown10.38 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD		R.T.
10.38	5.14 ng	308288	Acenaphthene-d10		9.79
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Metacetamol	151	C8H9NO2	000621-42-1	43
2	Pyridine, 2-butyl-, 1-oxide	151	C9H13NO	031396-32-4	38
3	m-Isopropoxvaniline	151	C9H13NO	041406-00-2	38
4	Acetaminophen	151	C8H9NO2	000103-90-2	38
5	p-Isopropoxyaniline	151	C9H13NO	007664-66-6	32



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF041020\
Data File : BF119910.D
Acq On : 10 Apr 2020 22:15
Operator : MA/SJ
Sample : L2199-05
Misc :
ALS Vial : 19 Sample Multiplier: 1

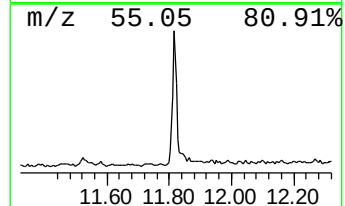
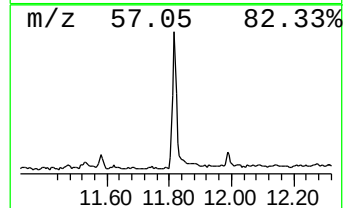
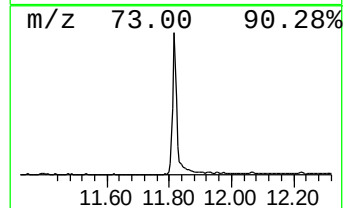
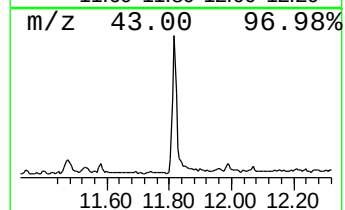
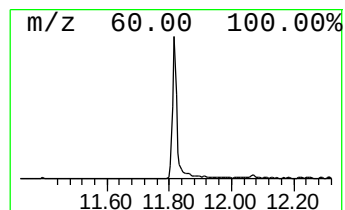
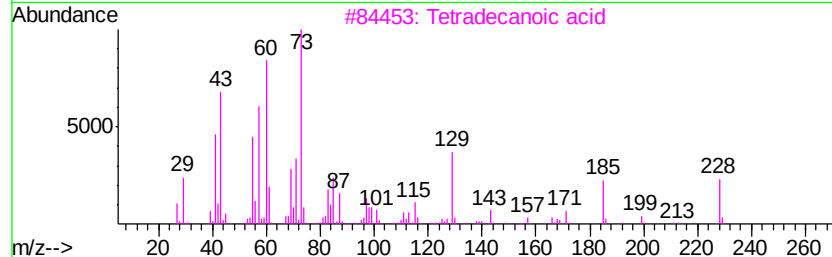
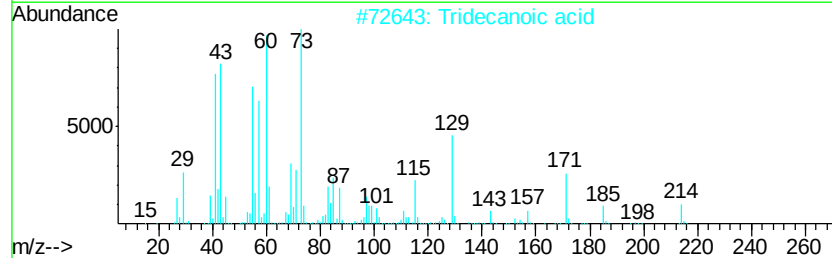
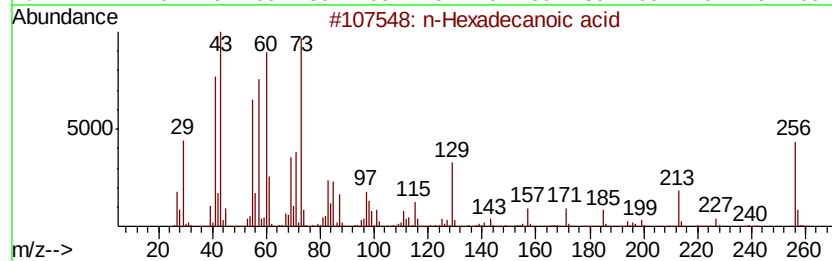
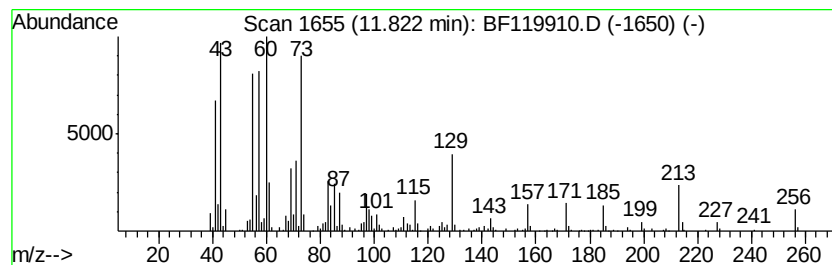
Instrument :
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ClientSampleId :
TP-39-38-COMP

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF040820.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD		R.T.
11.82	12.33 ng	665125	Phenanthrene-d10		11.28
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	94
2	Tridecanoic acid	214	C13H26O2	000638-53-9	94
3	Tetradecanoic acid	228	C14H28O2	000544-63-8	90
4	Pentadecanoic acid	242	C15H30O2	001002-84-2	83
5	Undecanoic acid	186	C11H22O2	000112-37-8	62



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF041020\
Data File : BF119910.D
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Operator : MA/SJ
Sample : L2199-05
Misc :
ALS Vial : 19 Sample Multiplier: 1

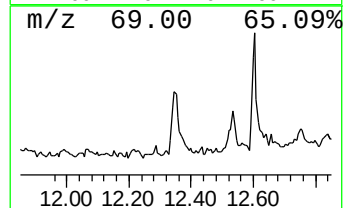
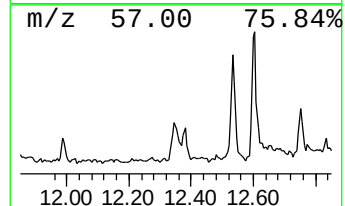
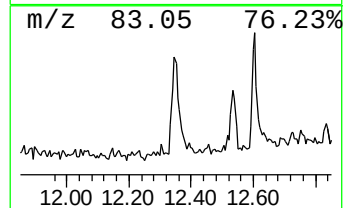
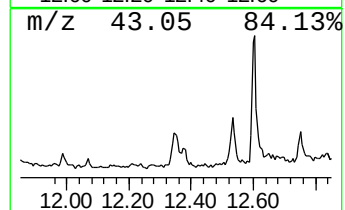
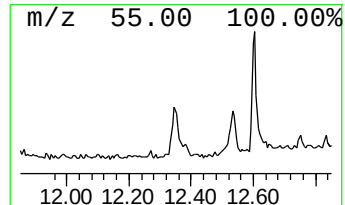
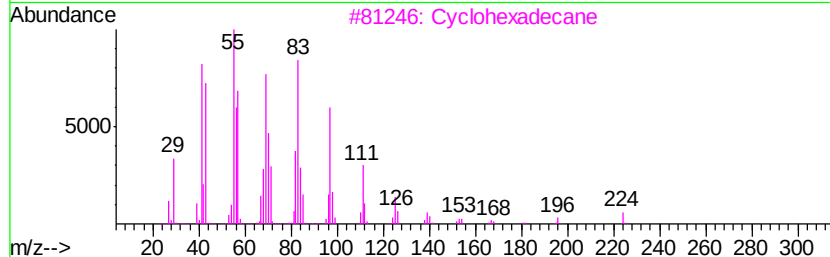
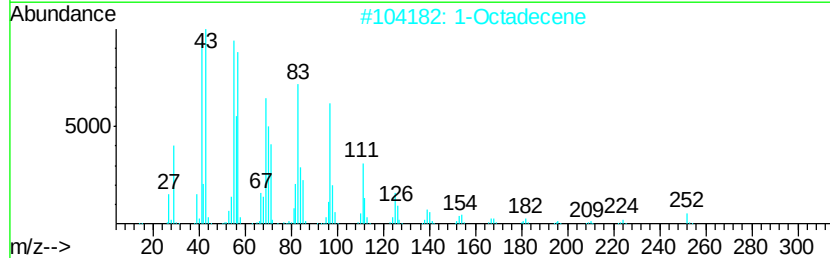
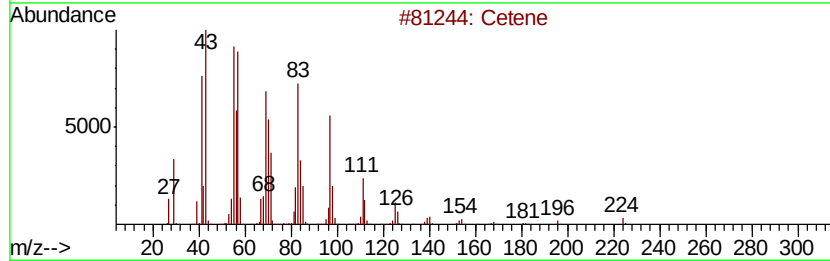
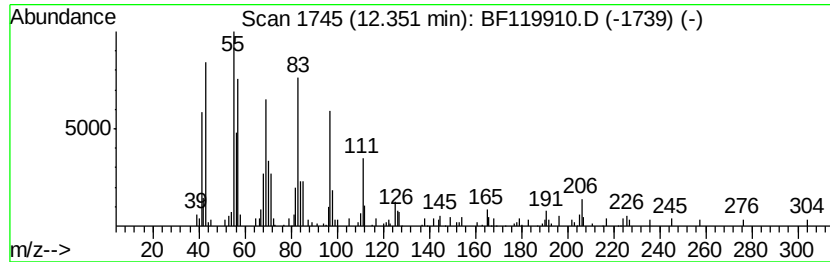
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF040820.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Cetene Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.35	4.13 ng	222962	Phenanthrene-d10	11.28
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Cetene	224 C16H32	000629-73-2 98
2		1-Octadecene	252 C18H36	000112-88-9 96
3		Cyclohexadecane	224 C16H32	000295-65-8 96
4		3-Octadecene, (E)-	252 C18H36	007206-19-1 95
5		Pentafluoropropionic acid, 4-hex...	388 C19H33F5O2	1000283-04-1 81



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF041020\
Data File : BF119910.D
Acq On : 10 Apr 2020 22:15
Operator : MA/SJ
Sample : L2199-05
Misc :
ALS Vial : 19 Sample Multiplier: 1

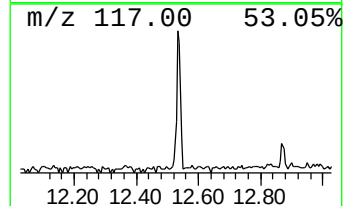
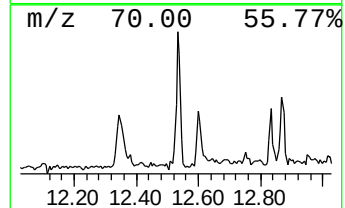
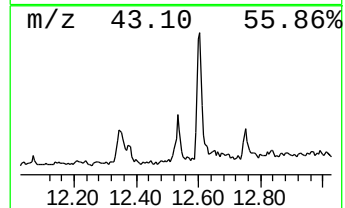
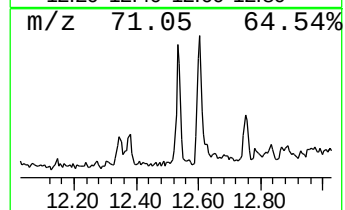
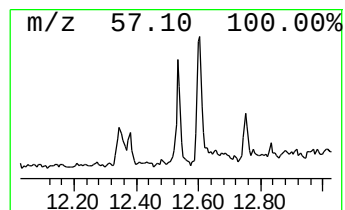
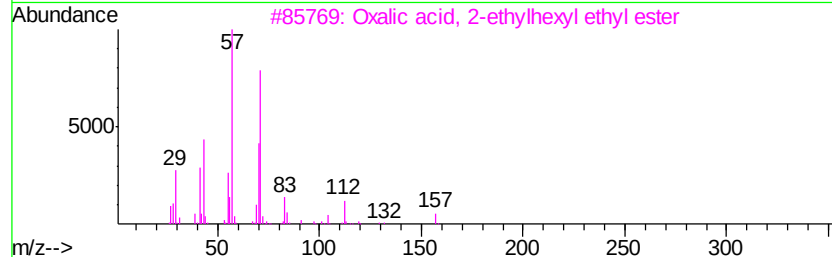
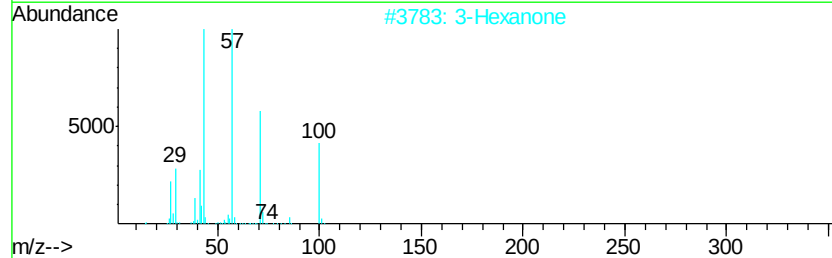
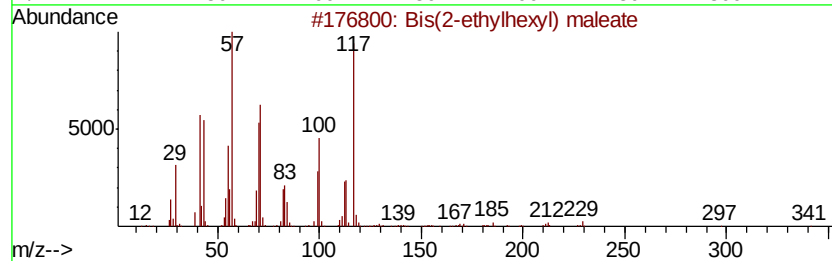
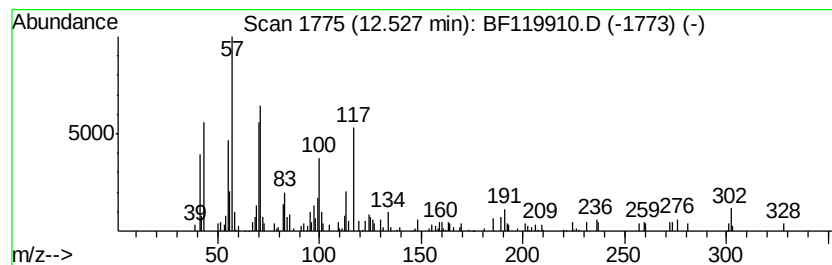
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ClientSampleId :
TP-39-38-COMP

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF040820.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 unknown12.53 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.53	3.00 ng	161665	Phenanthrene-d10	11.28	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Bis(2-ethylhexyl) maleate	340	C20H36O4	000142-16-5	38
2	3-Hexanone	100	C6H12O	000589-38-8	38
3	Oxalic acid, 2-ethylhexyl ethyl ...	230	C12H22O4	1000309-38-4	35
4	Sulfurous acid, decyl 2-ethylhex...	334	C18H38O3S	1000309-19-3	30
5	Sulfurous acid, 2-ethylhexyl tet...	390	C22H46O3S	1000309-19-7	30



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF041020\
Data File : BF119910.D
Acq On : 10 Apr 2020 22:15
Operator : MA/SJ
Sample : L2199-05
Misc :
ALS Vial : 19 Sample Multiplier: 1

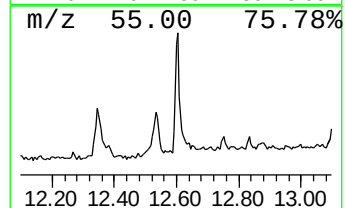
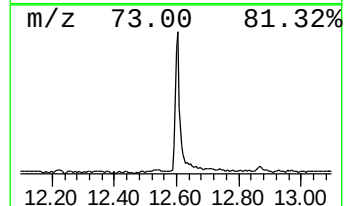
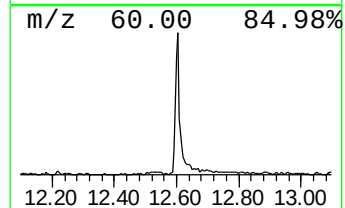
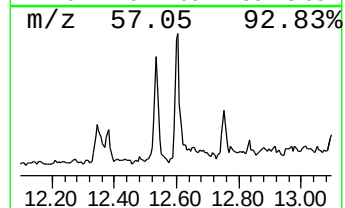
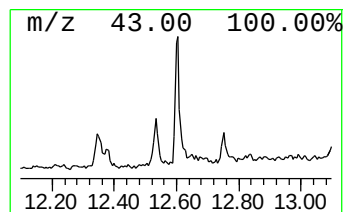
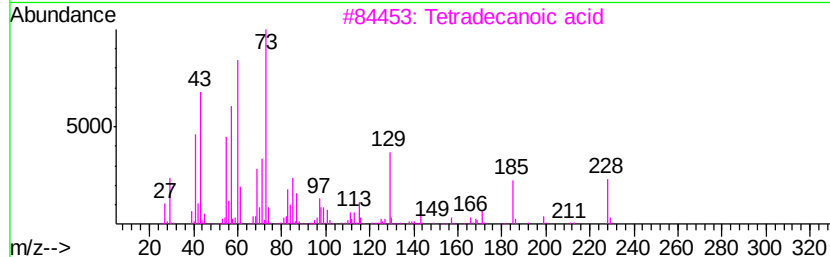
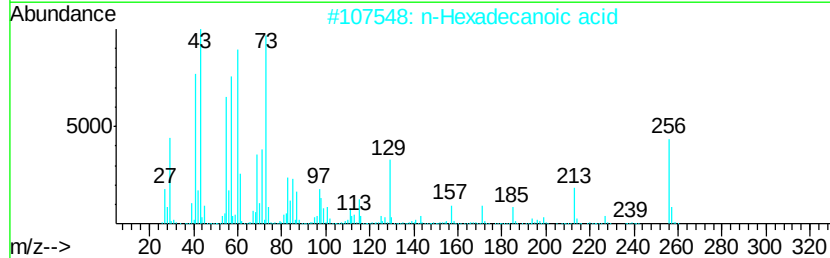
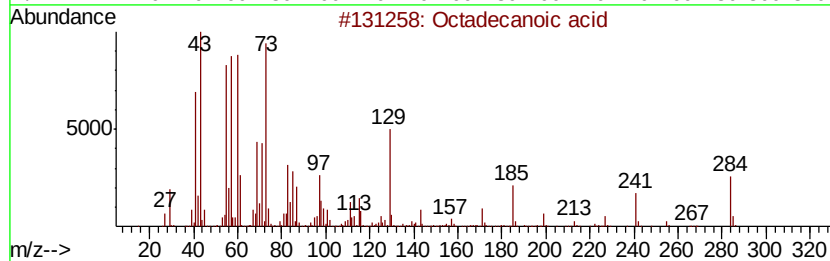
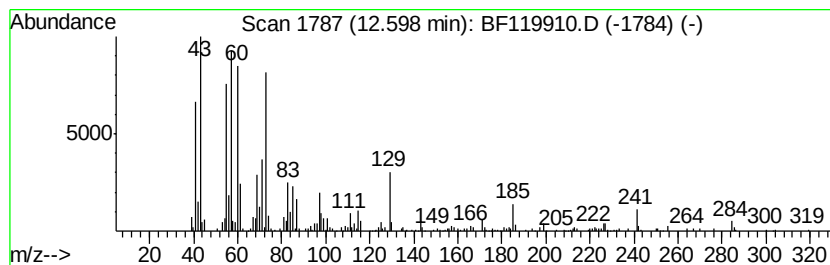
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF040820.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Octadecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.60	8.89 ng	473620	Chrysene-d12	13.92	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecanoic acid	284	C18H36O2	000057-11-4	99
2	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	90
3	Tetradecanoic acid	228	C14H28O2	000544-63-8	83
4	Pentadecanoic acid	242	C15H30O2	001002-84-2	74
5	n-Decanoic acid	172	C10H20O2	000334-48-5	70



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF041020\
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Sample : L2199-05
Misc :
ALS Vial : 19 Sample Multiplier: 1

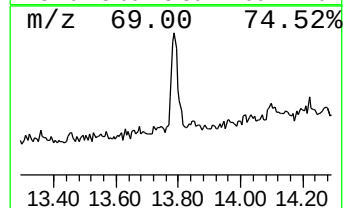
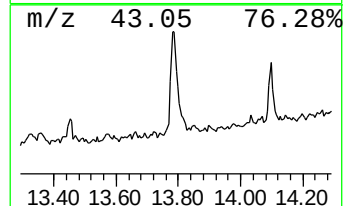
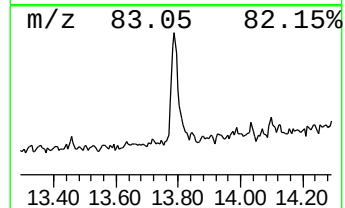
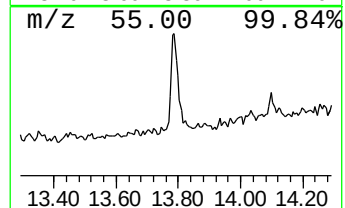
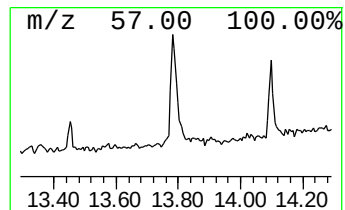
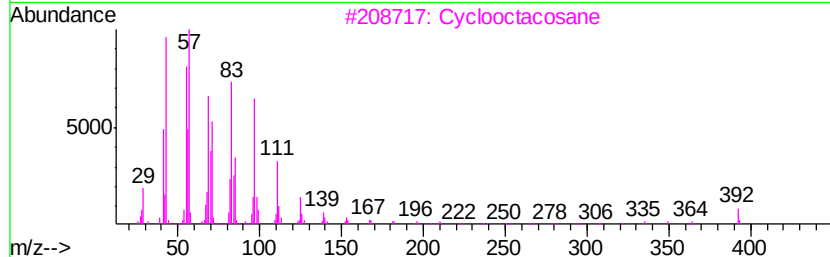
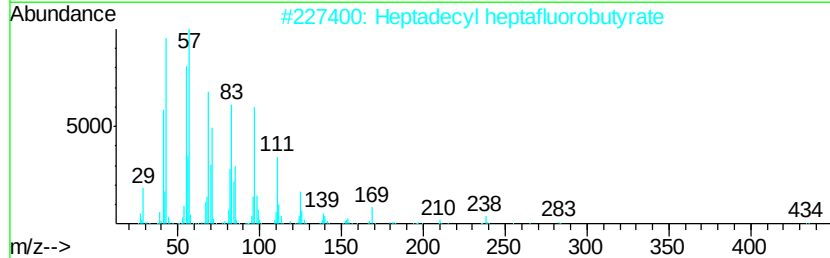
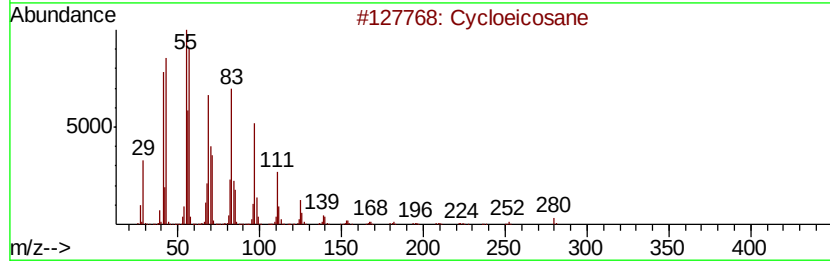
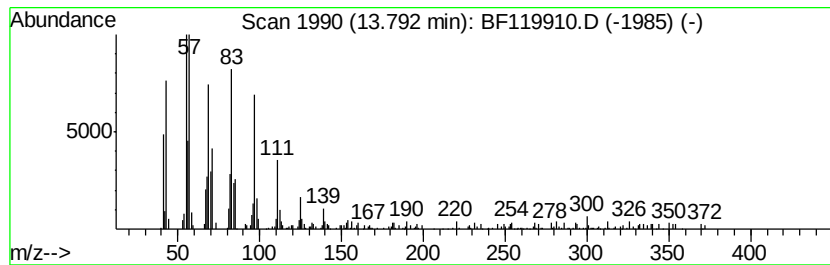
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF040820.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Cycloeicosane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.79	6.53 ng	348193	Chrysene-d12	13.92	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cvcloeicosane	280	C20H40	000296-56-0	93
2	Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	91
3	Cvclooctacosane	392	C28H56	000297-24-5	91
4	Heptadecyl trifluoroacetate	352	C19H35F3O2	1000351-87-0	91
5	1-Tricosene	322	C23H46	018835-32-0	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF041020\
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Sample : L2199-05
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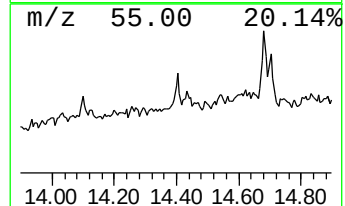
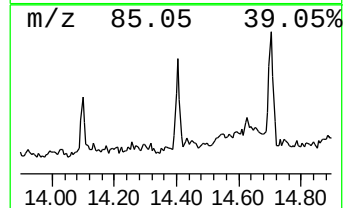
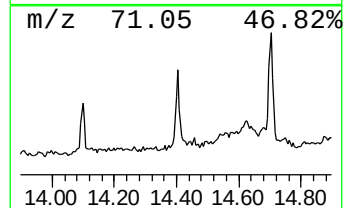
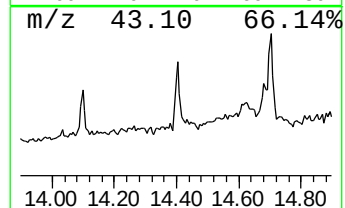
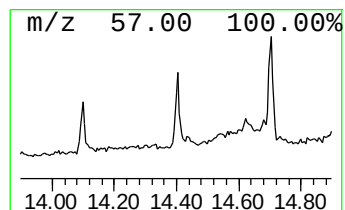
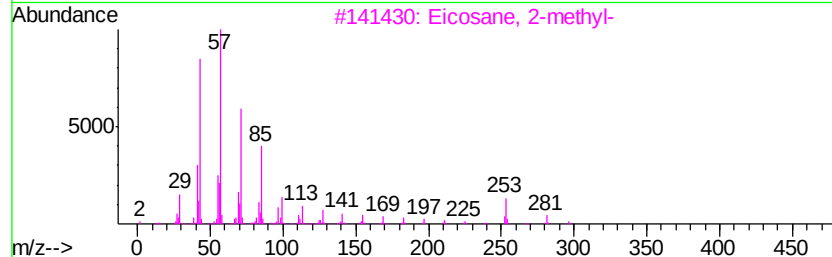
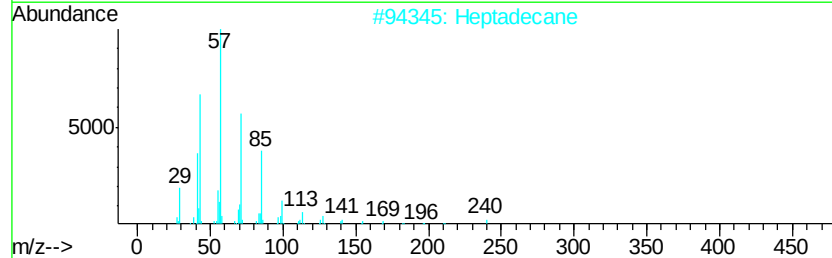
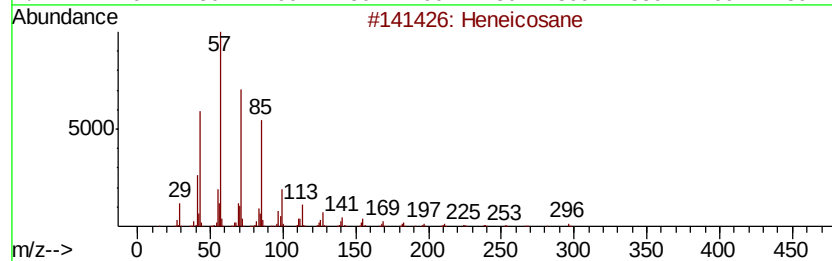
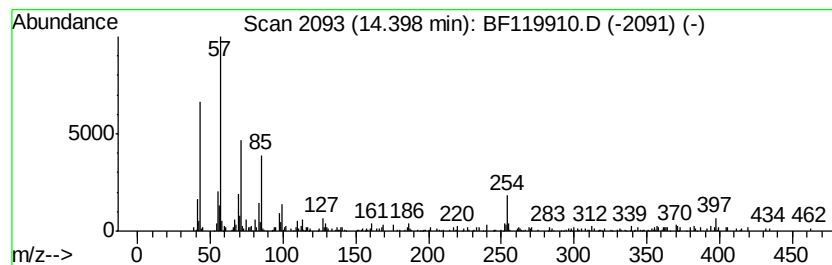
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Heneicosane Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.40	2.60 ng	138512	Chrysene-d12	13.92
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Heneicosane	296 C21H44	000629-94-7	95
2	Heptadecane	240 C17H36	000629-78-7	90
3	Eicosane, 2-methyl-	296 C21H44	001560-84-5	70
4	Pentacosane	352 C25H52	000629-99-2	70
5	3,5-Dimethyldodecane	198 C14H30	107770-99-0	68



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF041020\
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 Acq On : 10 Apr 2020 22:15
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 Misc :
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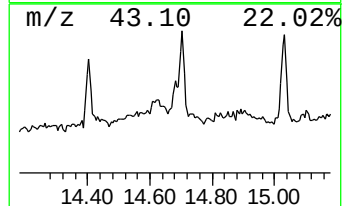
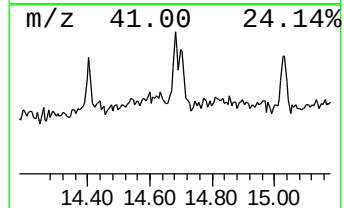
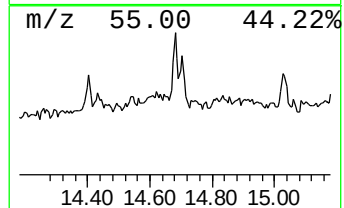
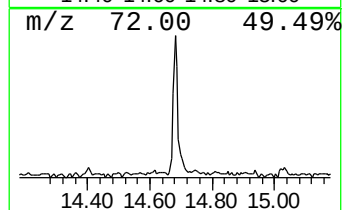
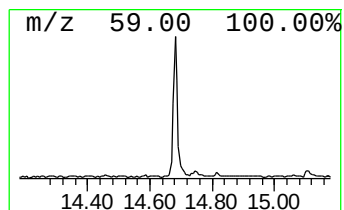
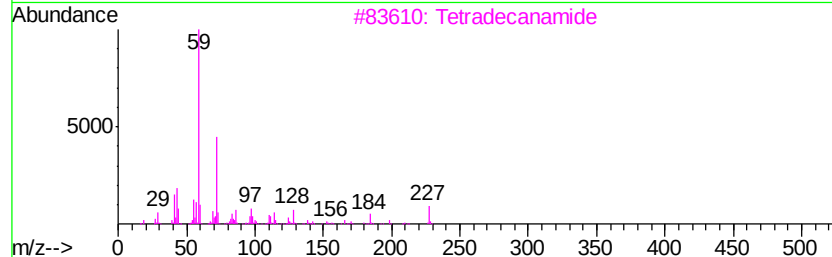
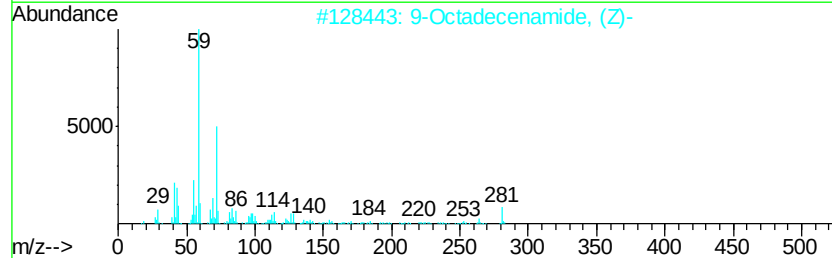
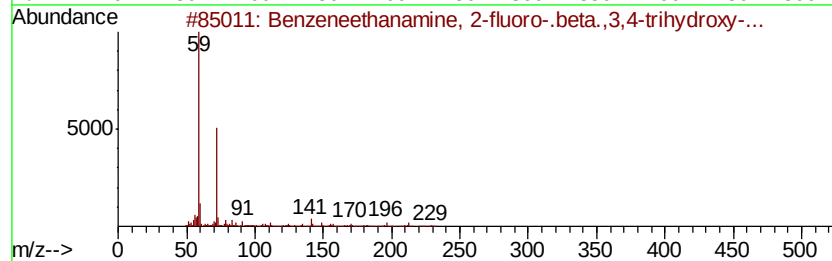
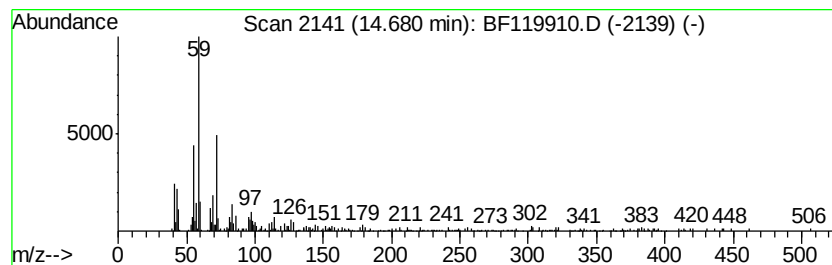
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

 Peak Number 14 Benzeneethanamine, 2-fluoro... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.68	4.52 ng	151164	Perylene-d12	15.36

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzeneethanamine, 2-fluoro-.bet...	229	C11H16FN03	061338-98-5	80
2		9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	80
3		Tetradecanamide	227	C14H29NO	000638-58-4	72
4		Hexadecanamide	255	C16H33NO	000629-54-9	72
5		Dodecanamide	199	C12H25NO	001120-16-7	58



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF041020\
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Acq On : 10 Apr 2020 22:15
Operator : MA/SJ
Sample : L2199-05
Misc :
ALS Vial : 19 Sample Multiplier: 1

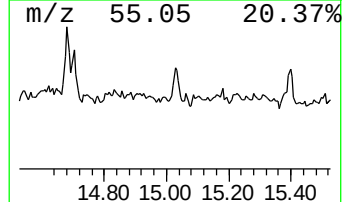
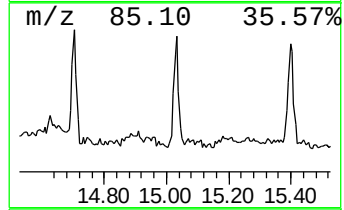
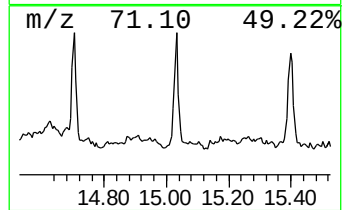
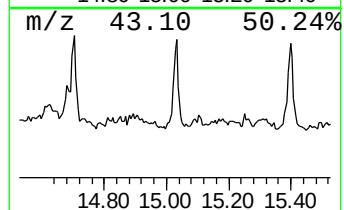
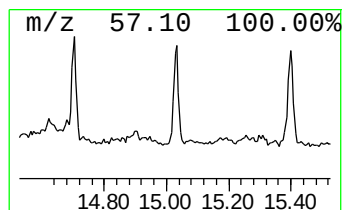
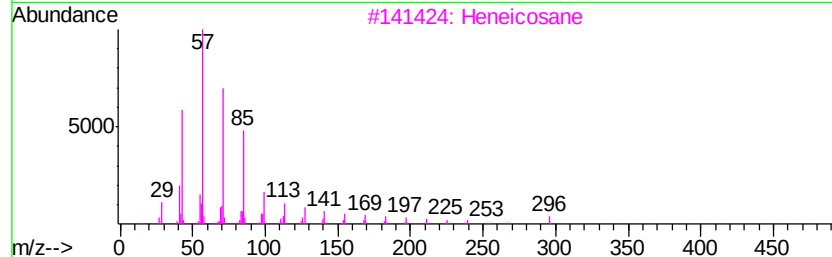
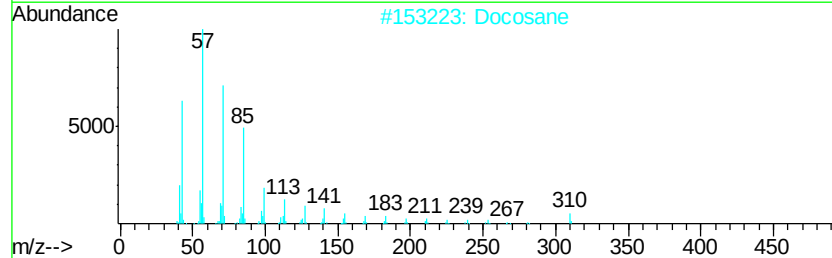
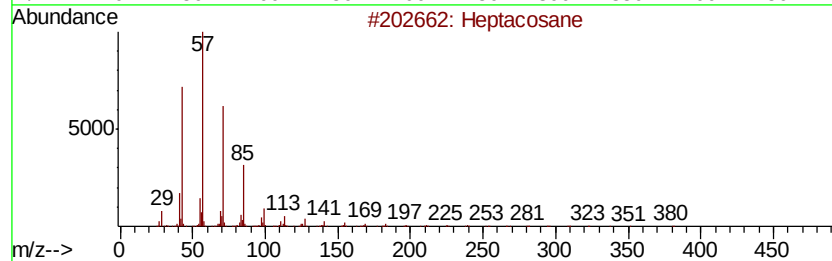
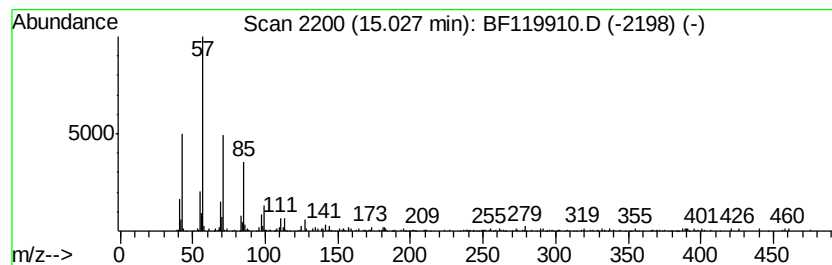
Instrument :
BNA_F
ClientSampleId :
TP-39-38-COMP

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF040820.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Heptacosane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
15.03	6.92 ng	231651	Perylene-d12	15.36	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptacosane	380	C27H56	000593-49-7	96
2	Docosane	310	C22H46	000629-97-0	94
3	Heneicosane	296	C21H44	000629-94-7	91
4	Tridecane, 1-iodo-	310	C13H27I	035599-77-0	87
5	Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	83



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF041020\
 Data File : BF119910.D
 Acq On : 10 Apr 2020 22:15
 Operator : MA/SJ
 Sample : L2199-05
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-39-38-COMP

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF040820.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
Propane, 2-methyl...	4.95	9.0	ng	465303	1	6.76	1034920	20.0
Ethanol, 2-(2-eth...	6.58	3.4	ng	177173	1	6.76	1034920	20.0
2,4,7,9-Tetrameth...	9.24	20.2	ng	1211310	3	9.79	1200100	20.0
1-Nonene	9.60	5.2	ng	311135	3	9.79	1200100	20.0
unknown10.38	10.38	5.1	ng	308288	3	9.79	1200100	20.0
n-Hexadecanoic acid	11.82	12.3	ng	665125	4	11.28	1078550	20.0
Cetene	12.35	4.1	ng	222962	4	11.28	1078550	20.0
unknown12.53	12.53	3.0	ng	161665	4	11.28	1078550	20.0
Octadecanoic acid	12.60	8.9	ng	473620	5	13.92	1065850	20.0
Cycloeicosane	13.79	6.5	ng	348193	5	13.92	1065850	20.0
Heneicosane	14.40	2.6	ng	138512	5	13.92	1065850	20.0
Benzeneethanamine...	14.68	4.5	ng	151164	6	15.36	669397	20.0
Heptacosane	15.03	6.9	ng	231651	6	15.36	669397	20.0