

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF052920\
 Data File : BF120447.D
 Acq On : 29 May 2020 18:55
 Operator : MA/SJ
 Sample : L2773-08
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 NM59-COMP-2020-0-6

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-BF052820.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	1.934	4	8	10	rBV	486949	511448	8.03%	0.886%
2	1.958	10	12	23	rVB	1831320	2154208	33.82%	3.734%
3	2.110	33	38	54	rVB	1146200	1475073	23.16%	2.557%
4	2.228	54	58	75	rVB5	45119	131171	2.06%	0.227%
5	3.310	239	242	256	rBV	27571	81716	1.28%	0.142%
6	5.457	601	607	610	rBV	4252348	4699432	73.78%	8.145%
7	6.463	773	778	781	rBV	4498023	5091241	79.93%	8.824%
8	6.663	809	812	827	rBV	716694	602537	9.46%	1.044%
9	6.840	838	842	845	rBV	1834806	1671580	26.24%	2.897%
10	6.992	864	868	872	rVB	4798795	4527159	71.07%	7.847%
11	7.398	931	937	940	rBV	3601026	3613097	56.72%	6.262%
12	8.116	1055	1059	1063	rBV	2506363	2099855	32.97%	3.640%
13	9.198	1237	1243	1246	rBV	6562236	6369782	100.00%	11.040%
14	9.310	1254	1262	1266	rBV	1716265	1692389	26.57%	2.933%
15	9.586	1305	1309	1312	rBV	1112769	845826	13.28%	1.466%
16	9.875	1353	1358	1361	rBV	2477132	2277004	35.75%	3.947%
17	10.457	1453	1457	1464	rBV	140007	160061	2.51%	0.277%
18	10.657	1486	1491	1495	rBV	3696620	3533091	55.47%	6.124%
19	11.357	1605	1610	1612	rBV	2339980	1996430	31.34%	3.460%
20	11.380	1612	1614	1617	rVV	293500	230585	3.62%	0.400%
21	11.428	1617	1622	1626	rVB	90130	93483	1.47%	0.162%
22	11.857	1691	1695	1697	rBV2	70664	75678	1.19%	0.131%
23	11.886	1697	1700	1703	rVV	175157	159882	2.51%	0.277%
24	11.969	1710	1714	1716	rBV2	83714	83669	1.31%	0.145%
25	12.416	1785	1790	1793	rBV2	85002	142515	2.24%	0.247%
26	12.445	1793	1795	1799	rVB2	67274	73835	1.16%	0.128%
27	12.563	1810	1815	1819	rBV	562222	586040	9.20%	1.016%
28	12.604	1819	1822	1825	rBV2	109406	87404	1.37%	0.151%
29	12.792	1848	1854	1857	rBV	503016	511768	8.03%	0.887%
30	12.945	1875	1880	1883	rBV	5266758	4770351	74.89%	8.268%
31	13.121	1907	1910	1913	rVB	83014	79506	1.25%	0.138%
32	13.186	1914	1921	1924	rBV3	73988	150694	2.37%	0.261%
33	13.739	2010	2015	2018	rBV2	46201	75900	1.19%	0.132%
34	13.851	2028	2034	2040	rBV3	417844	709371	11.14%	1.230%

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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	13.992	2050	2058	2061	rBV	1760868	1856798	29.15%	3.218%
36	14.016	2061	2062	2068	rVB	283191	187915	2.95%	0.326%
37	14.480	2133	2141	2146	rBV4	132561	317907	4.99%	0.551%
38	14.916	2212	2215	2221	rVB3	60330	77455	1.22%	0.134%
39	15.021	2228	2233	2236	rBV	238421	361953	5.68%	0.627%
40	15.110	2244	2248	2250	rBV2	131341	166804	2.62%	0.289%
41	15.139	2250	2253	2258	rVB3	85279	144670	2.27%	0.251%
42	15.315	2280	2283	2289	rVB	145384	177157	2.78%	0.307%
43	15.380	2289	2294	2299	rBV	185054	229606	3.60%	0.398%
44	15.445	2300	2305	2309	rBV	1294484	1543156	24.23%	2.675%
45	15.915	2380	2385	2390	rBV	164248	249574	3.92%	0.433%
46	15.974	2390	2395	2408	rVB4	116736	295412	4.64%	0.512%
47	16.704	2514	2519	2527	rVB4	82704	154433	2.42%	0.268%
48	16.874	2543	2548	2552	rBV2	76595	145334	2.28%	0.252%
49	17.304	2617	2621	2629	rVB2	61667	109870	1.72%	0.190%
50	17.486	2645	2652	2665	rVB6	96617	313335	4.92%	0.543%

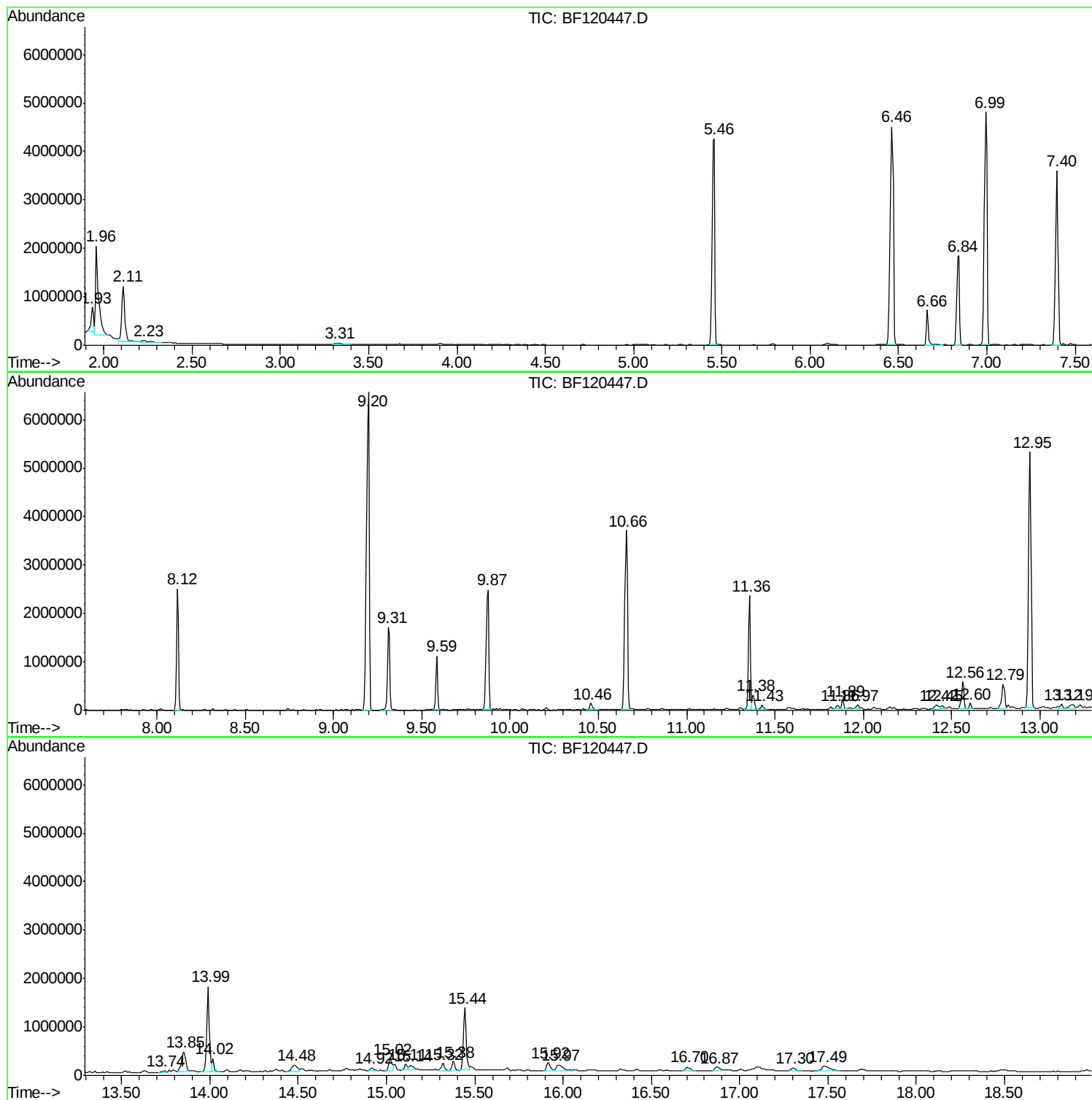
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF052820.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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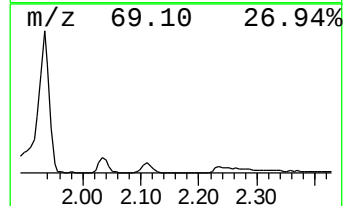
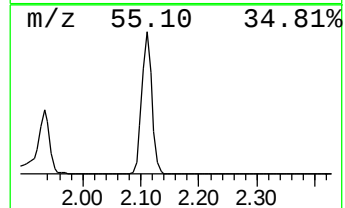
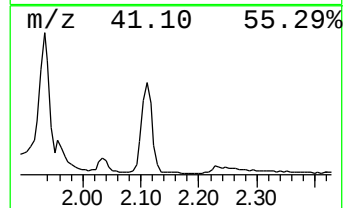
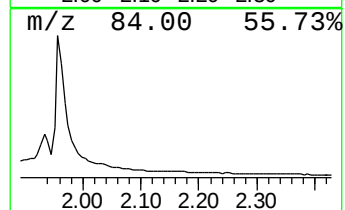
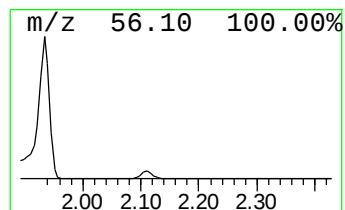
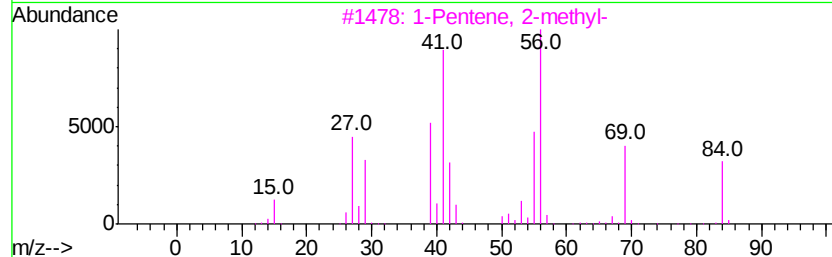
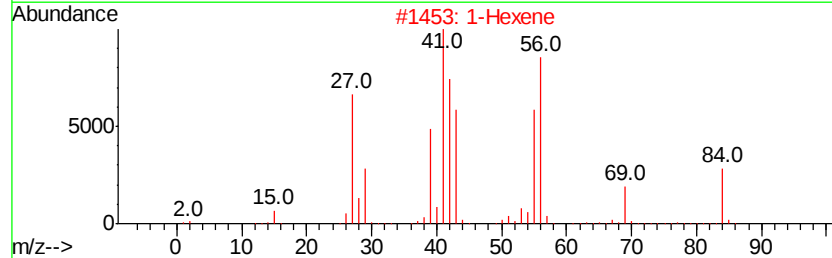
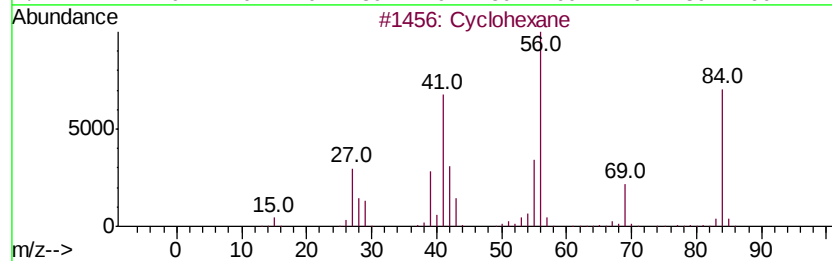
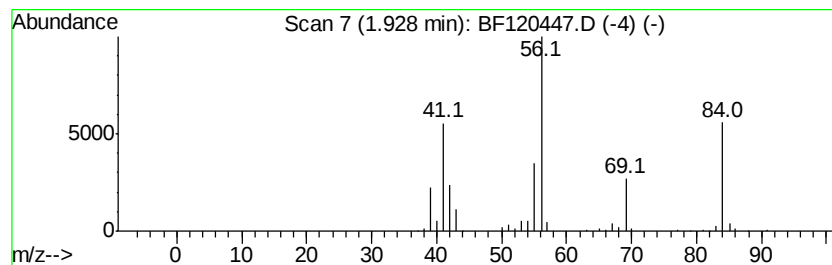
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF052820.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Cyclohexane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.93	6.12 ng	511448	1,4-Dichlorobenzene-d4	6.84

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexane	84	C6H12	000110-82-7	91
2		1-Hexene	84	C6H12	000592-41-6	59
3		1-Pentene, 2-methyl-	84	C6H12	000763-29-1	56
4		Cyclopentane, methyl-	84	C6H12	000096-37-7	50
5		2,2-Dimethylglutaric anhydride	142	C7H10O3	002938-48-9	40



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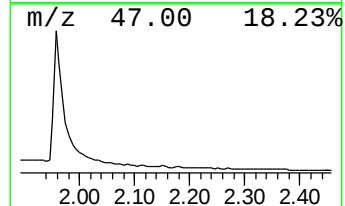
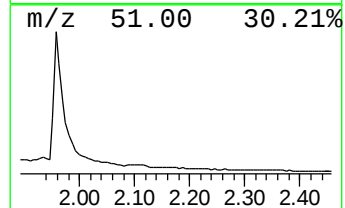
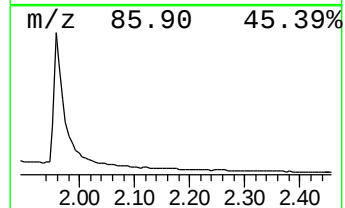
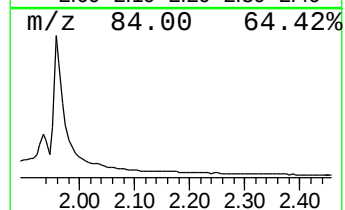
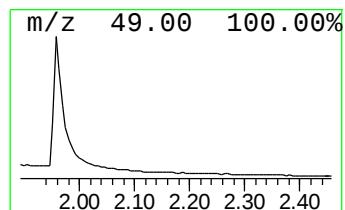
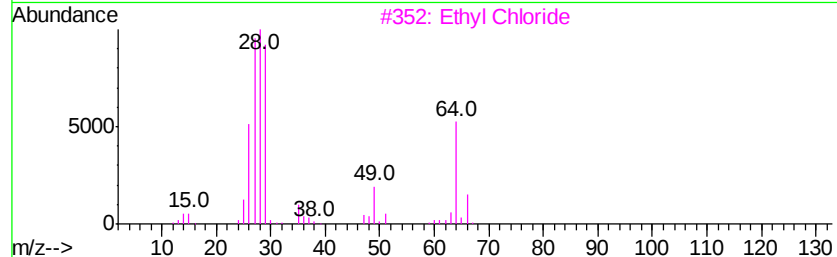
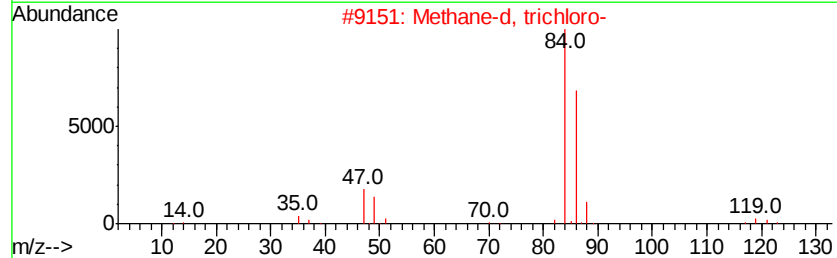
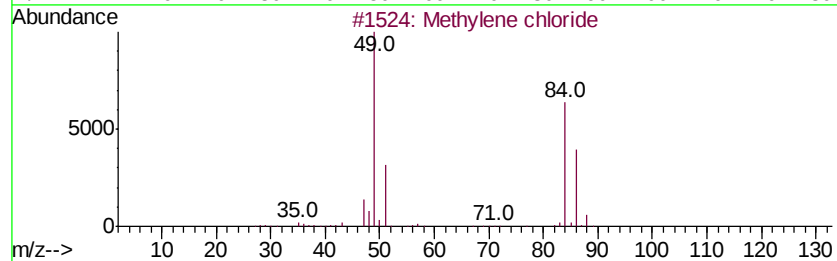
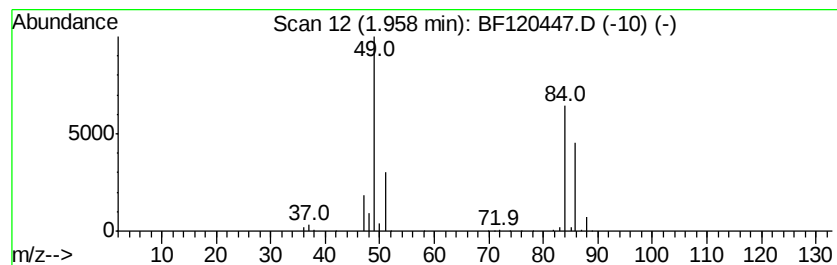
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 Methylene chloride Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.96	25.77 ng	2154210	1,4-Dichlorobenzene-d4	6.84

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Methylene chloride	84	CH2Cl2	000075-09-2	95
2		Methane-d, trichloro-	119	CDCl3	000865-49-6	9
3		Ethyl Chloride	64	C2H5Cl	000075-00-3	4
4		Acetic acid, chloro-, ethyl ester	122	C4H7ClO2	000105-39-5	2
5		2-Chloroethanol	80	C2H5ClO	000107-07-3	1



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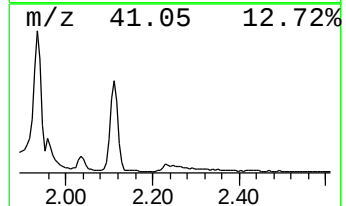
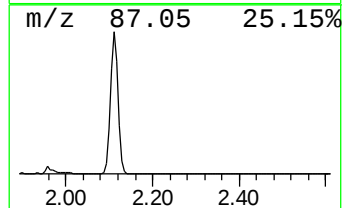
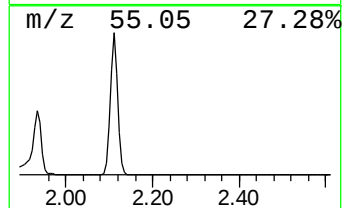
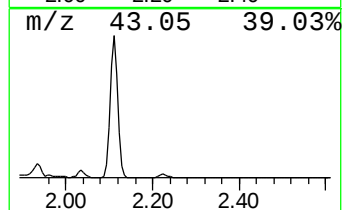
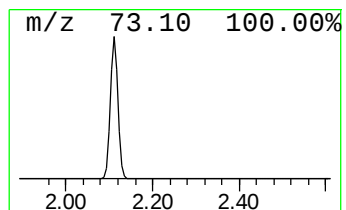
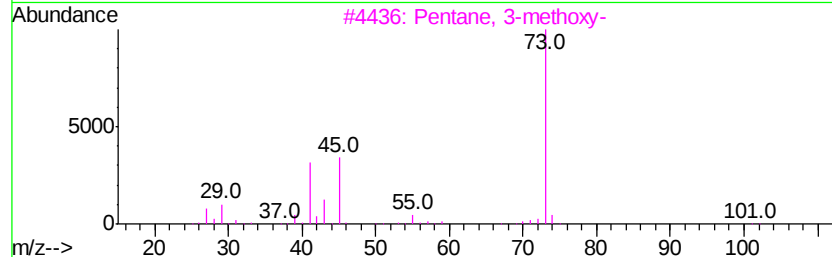
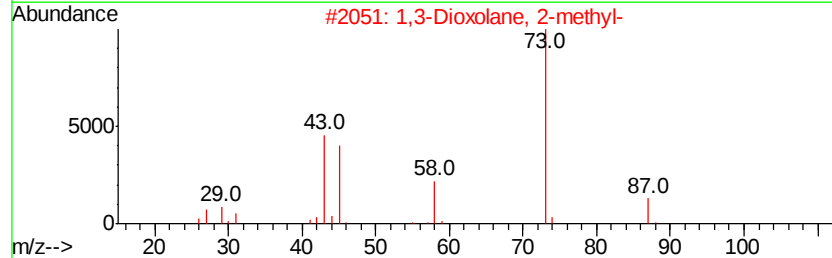
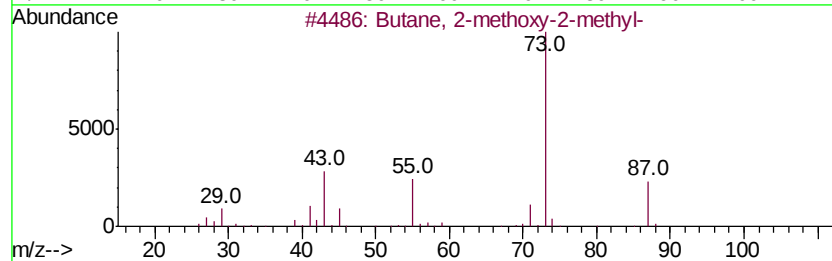
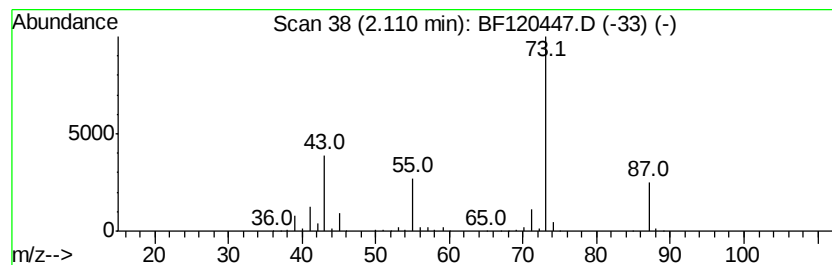
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Butane, 2-methoxy-2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.11	17.65 ng	1475070	1,4-Dichlorobenzene-d4	6.84

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2		1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	25
3		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	12
4		Silane, tetramethyl-	88	C4H12Si	000075-76-3	10
5		N-Ethylformamide	73	C3H7NO	000627-45-2	9



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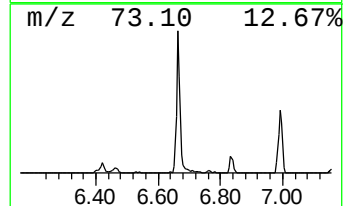
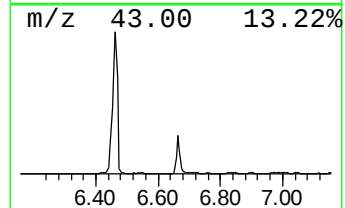
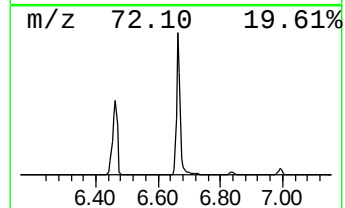
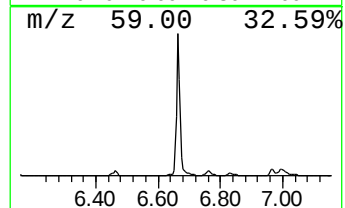
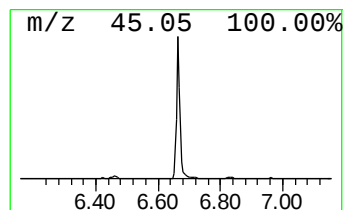
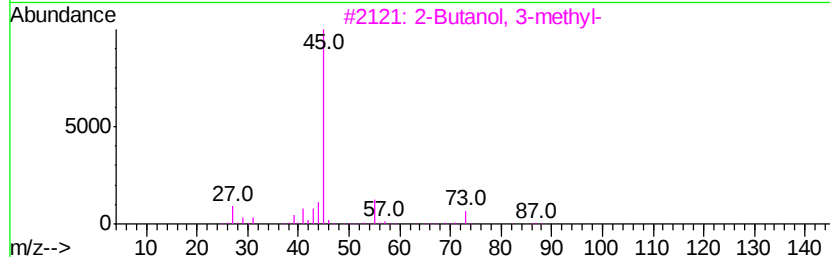
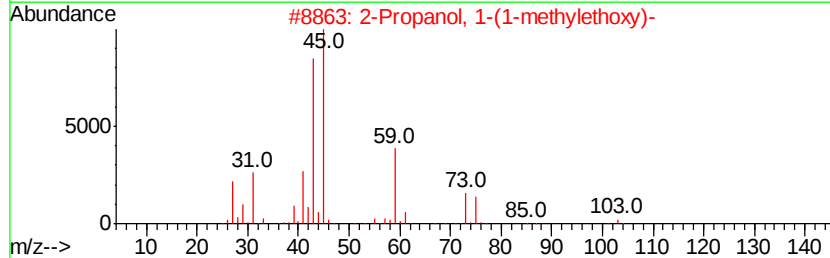
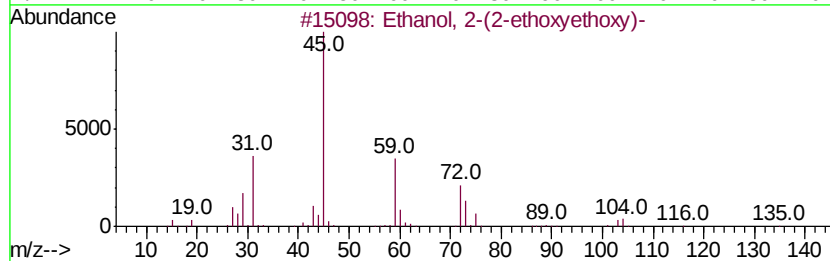
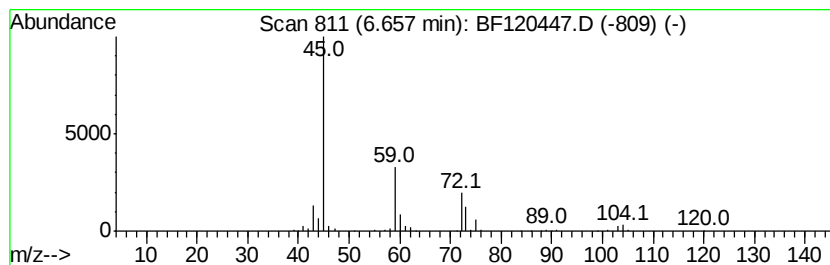
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Ethanol, 2-(2-ethoxyethoxy)- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
6.66	7.21 ng	602537	1,4-Dichlorobenzene-d4	6.84	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethanol, 2-(2-ethoxyethoxy)-	134	C6H14O3	000111-90-0	90
2	2-Propanol, 1-(1-methylethoxy)-	118	C6H14O2	003944-36-3	45
3	2-Butanol, 3-methyl-	88	C5H12O	000598-75-4	9
4	Acetic acid, hydroxy-, ethyl ester	104	C4H8O3	000623-50-7	9
5	Propanoic acid, 2-hydroxy-, meth...	104	C4H8O3	002155-30-8	9



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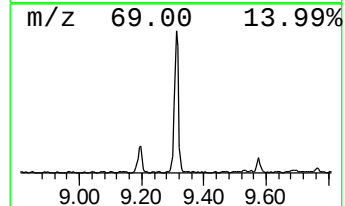
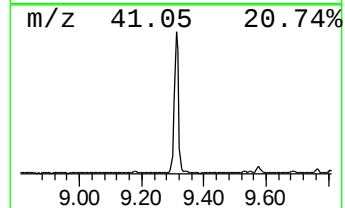
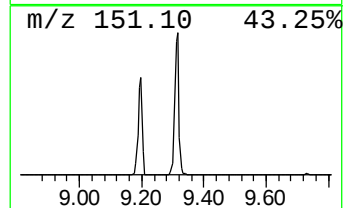
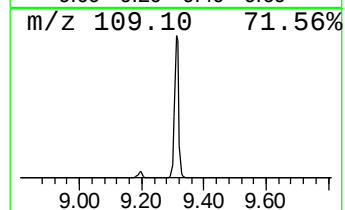
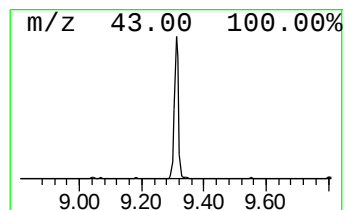
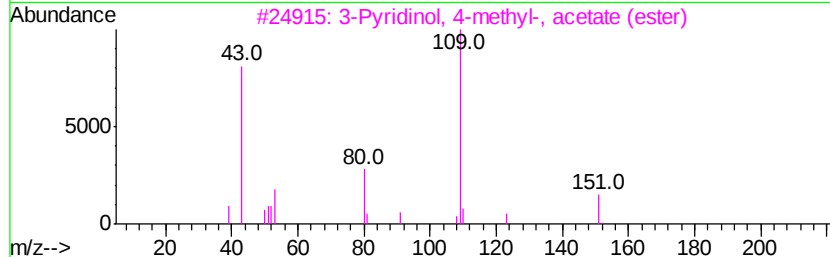
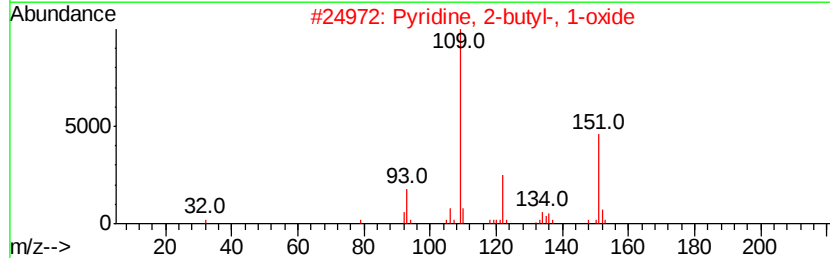
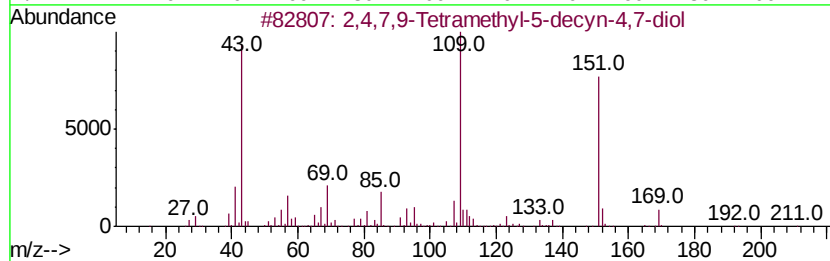
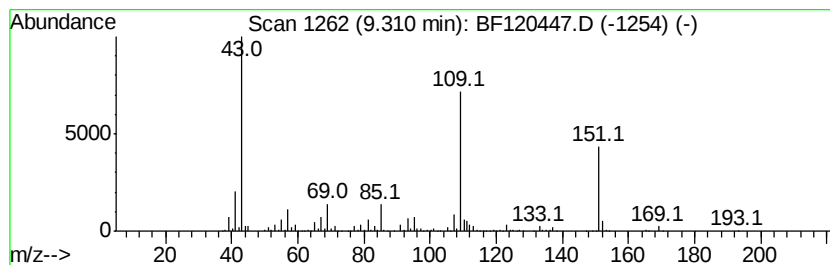
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 2,4,7,9-Tetramethyl-5-decyn... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.31	14.87 ng	1692390	Acenaphthene-d10	9.87

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2,4,7,9-Tetramethyl-5-decyn-4,7-...	226	C14H26O2	000126-86-3	90
2		Pyridine, 2-butyl-, 1-oxide	151	C9H13NO	031396-32-4	59
3		3-Pyridinol, 4-methyl-, acetate ...	151	C8H9NO2	001006-96-8	50
4		Metacetamol	151	C8H9NO2	000621-42-1	50
5		m-Isopropoxyaniline	151	C9H13NO	041406-00-2	50



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF052920\
Data File : BF120447.D
Acq On : 29 May 2020 18:55
Operator : MA/SJ
Sample : L2773-08
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
NM59-COMP-2020-0-6

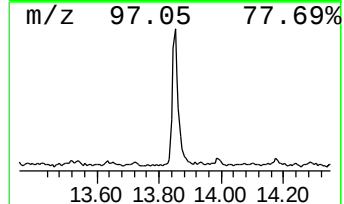
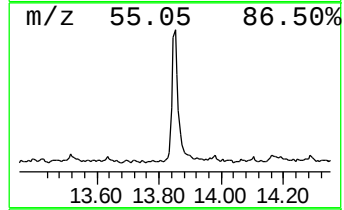
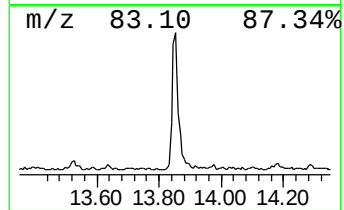
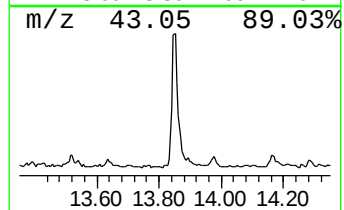
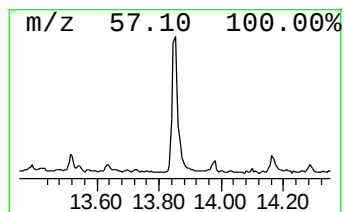
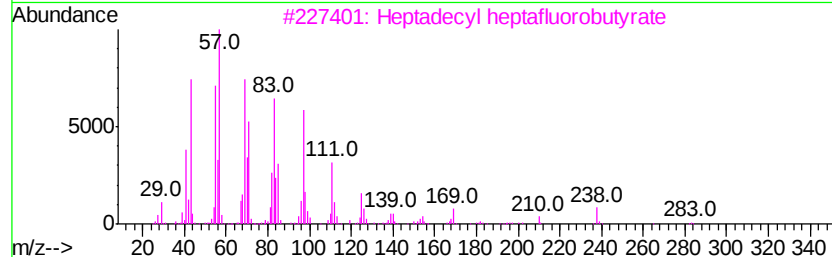
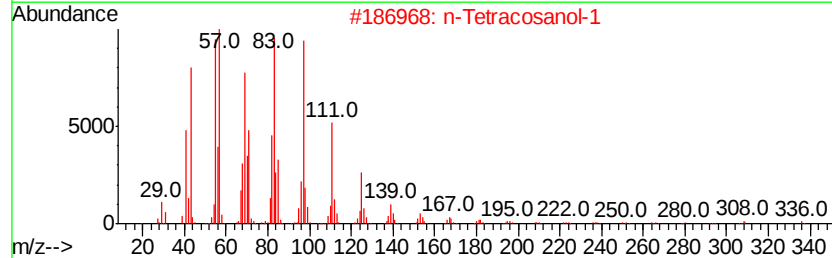
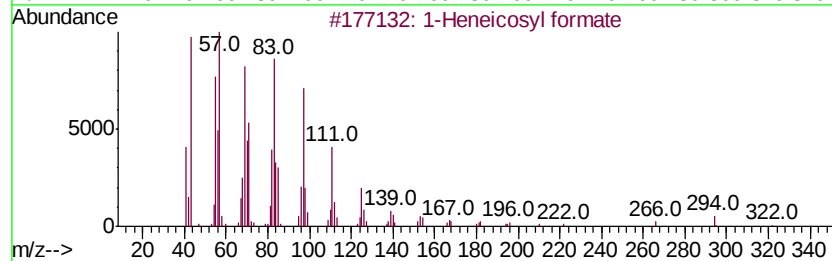
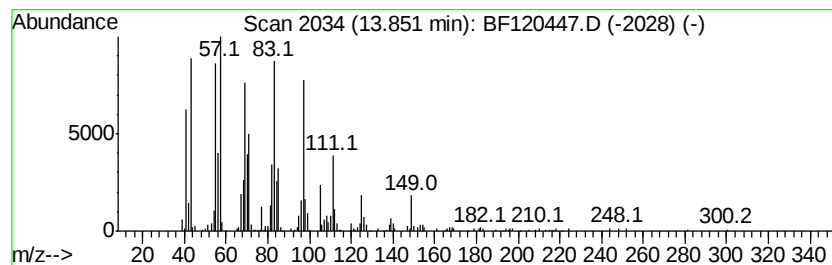
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 1-Heneicosyl formate Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.85	7.64 ng	709371	Chrysene-d12	13.99

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Heneicosyl formate	340	C22H44O2	077899-03-7	95
2		n-Tetracosanol-1	354	C24H50O	000506-51-4	94
3		Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	94
4		Trifluoroacetoxy hexadecane	338	C18H33F3O2	006222-03-3	94
5		Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	94



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF052920\
Data File : BF120447.D
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Misc :
ALS Vial : 16 Sample Multiplier: 1

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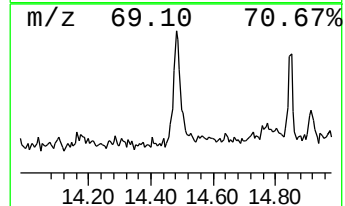
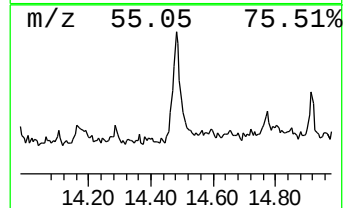
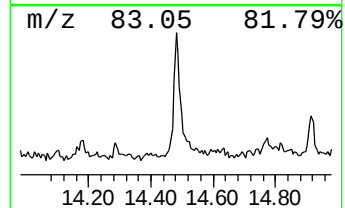
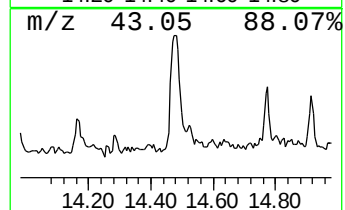
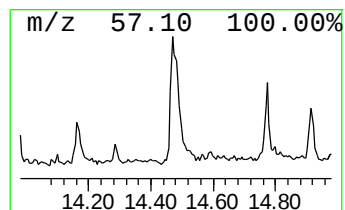
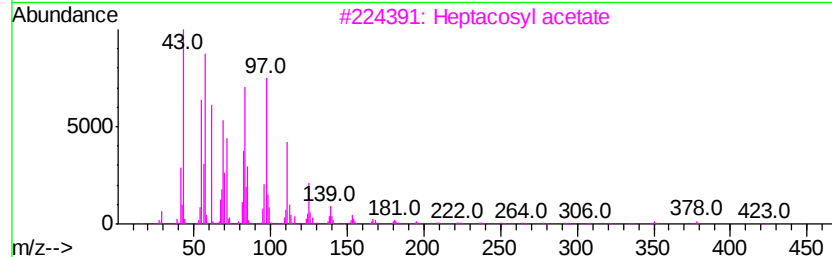
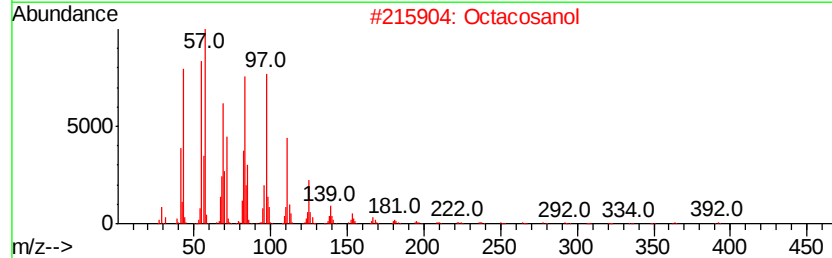
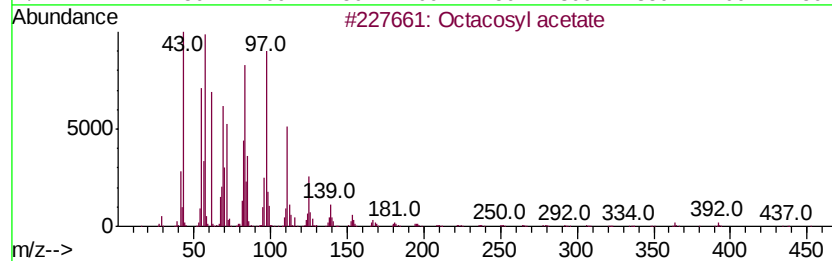
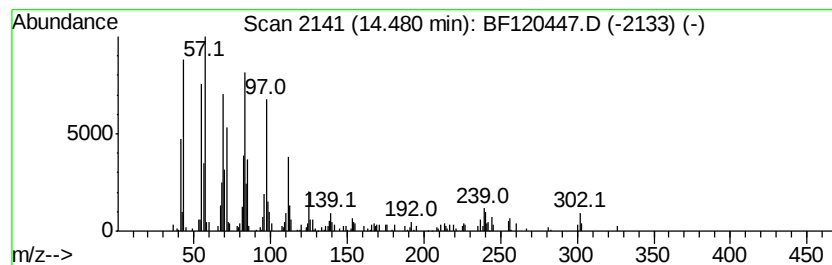
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Octacosyl acetate Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.48	3.42 ng	317907	Chrysene-d12	13.99

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Octacosyl acetate	452	C30H60O2	018206-97-8	93
2		Octacosanol	410	C28H58O	000557-61-9	93
3		Heptacosyl acetate	438	C29H58O2	1000351-78-2	93
4		1-Heptacosanol	396	C27H56O	002004-39-9	93
5		triacontyl acetate	480	C32H64O2	041755-58-2	93



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF052920\
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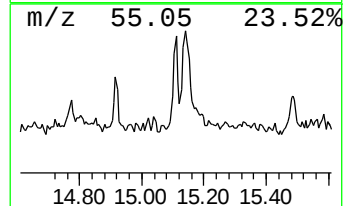
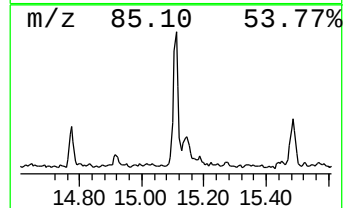
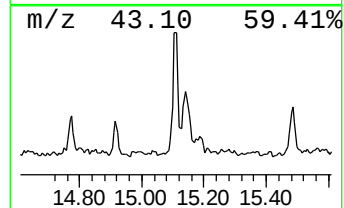
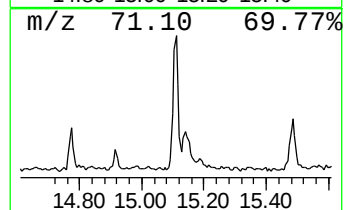
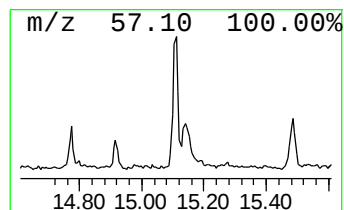
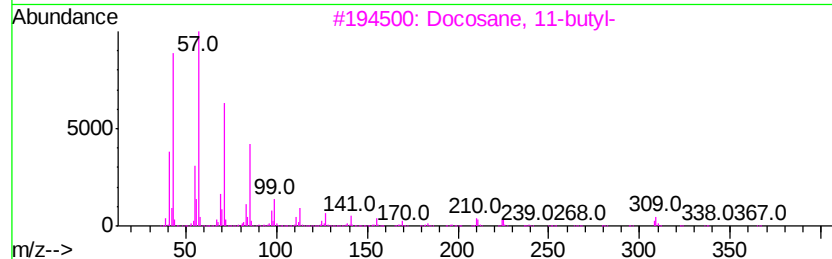
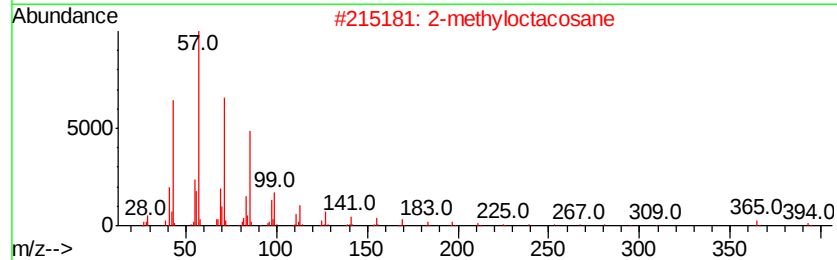
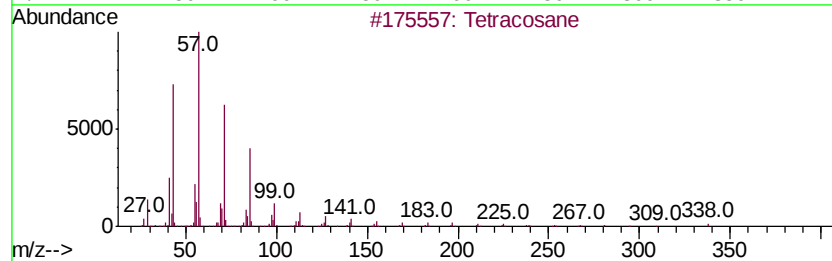
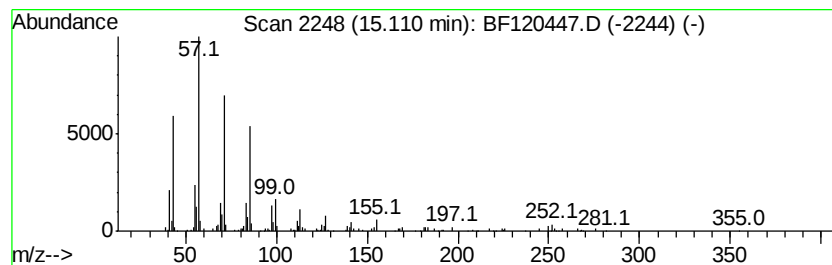
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Tetracosane Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.11	2.16 ng	166804	Perylene-d12	15.44

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetracosane	338	C24H50	000646-31-1	87
2		2-methyloctacosane	408	C29H60	1000376-72-8	87
3		Docosane, 11-butyl-	366	C26H54	013475-76-8	87
4		Heneicosane	296	C21H44	000629-94-7	87
5		Pentacosane	352	C25H52	000629-99-2	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF052920\
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Misc :
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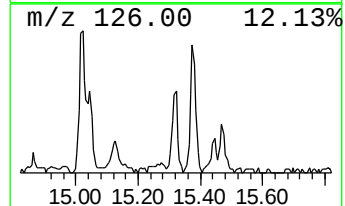
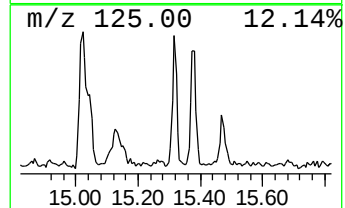
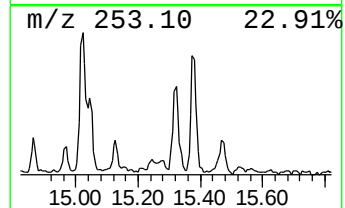
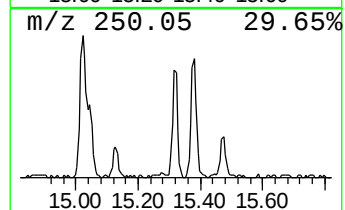
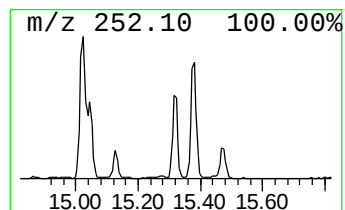
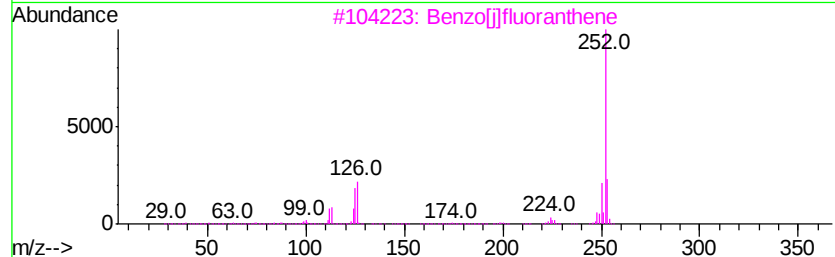
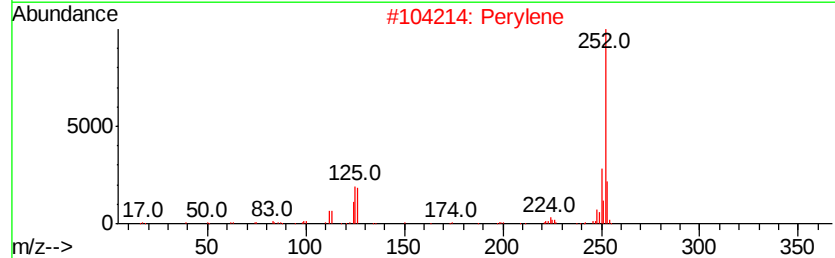
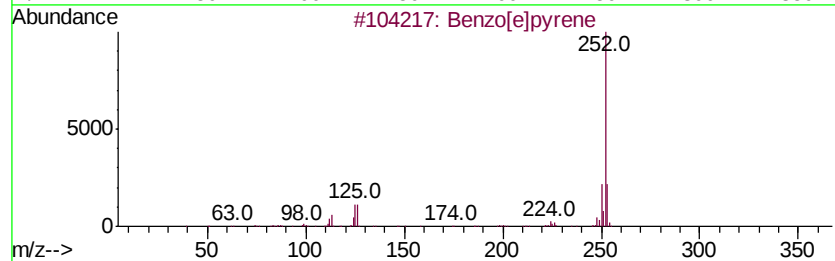
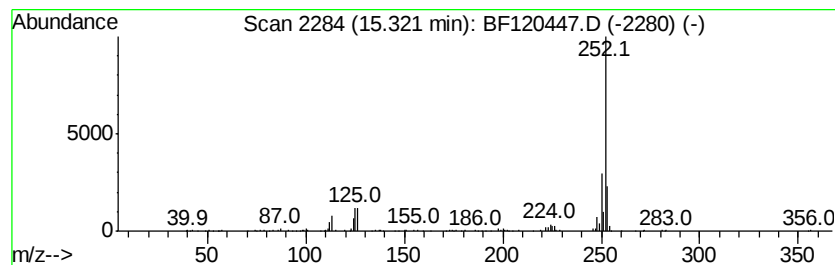
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Benzo[e]pyrene Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.	
15.32	2.30 ng	177157	Perylene-d12	15.44	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzo[e]pyrene	252	C20H12	000192-97-2	97
2	Perylene	252	C20H12	000198-55-0	95
3	Benzo[i]fluoranthene	252	C20H12	000205-82-3	93
4	Benzo[k]fluoranthene	252	C20H12	000207-08-9	91
5	Benz[e]acephenanthrylene	252	C20H12	000205-99-2	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF052920\
Data File : BF120447.D
Acq On : 29 May 2020 18:55
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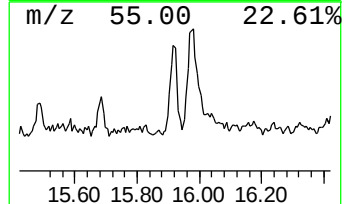
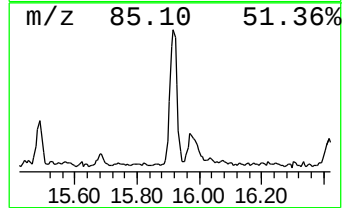
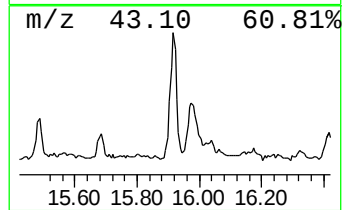
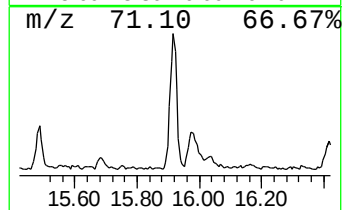
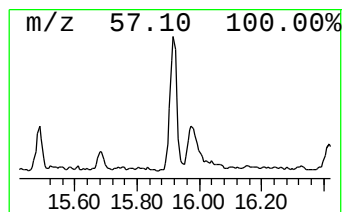
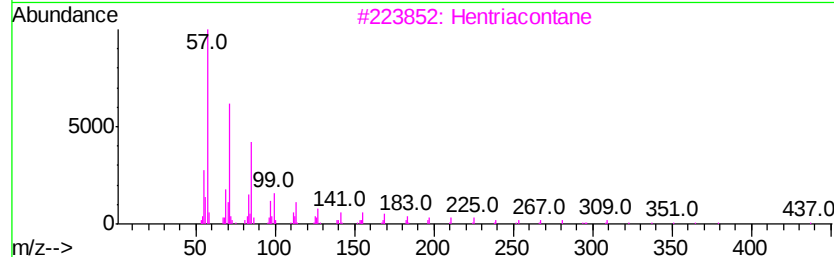
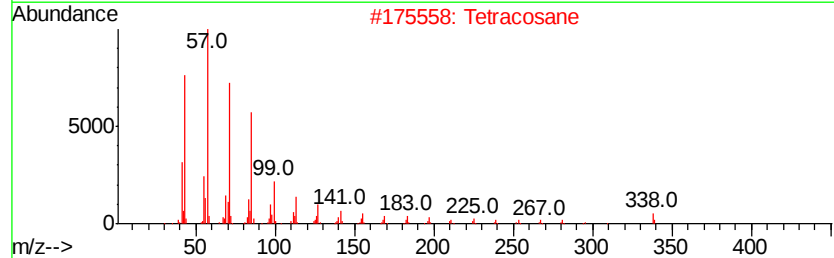
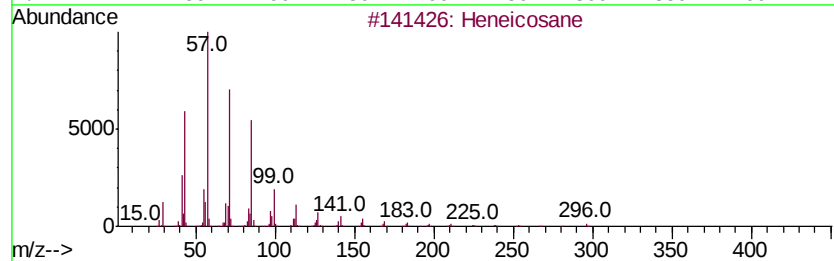
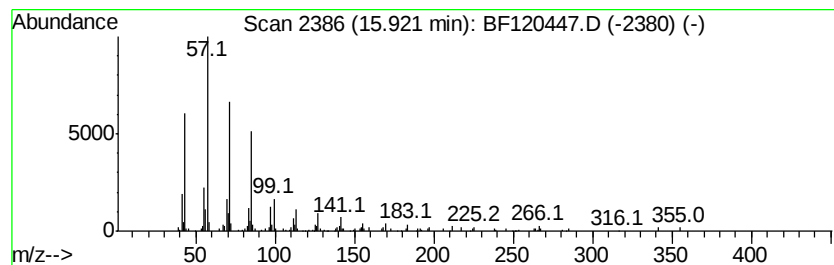
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Heneicosane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.92	3.23 ng	249574	Perylene-d12	15.44

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heneicosane	296	C21H44	000629-94-7	90
2			Tetracosane	338	C24H50	000646-31-1	90
3			Hentriacontane	437	C31H64	000630-04-6	90
4			Hexacosane	366	C26H54	000630-01-3	90
5			Octacosane	394	C28H58	000630-02-4	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF052920\
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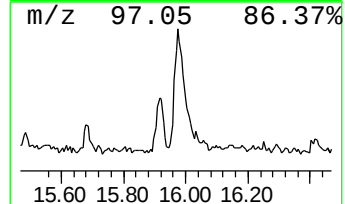
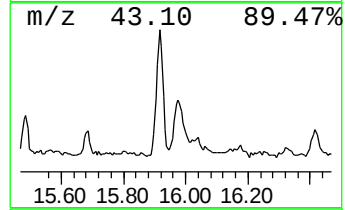
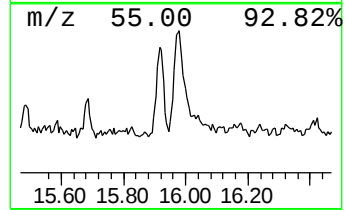
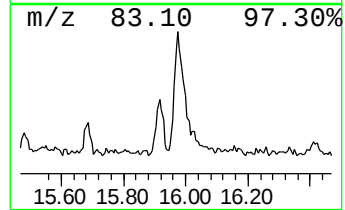
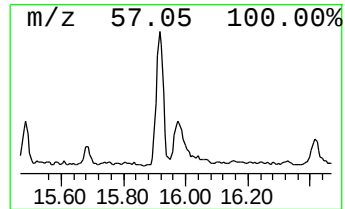
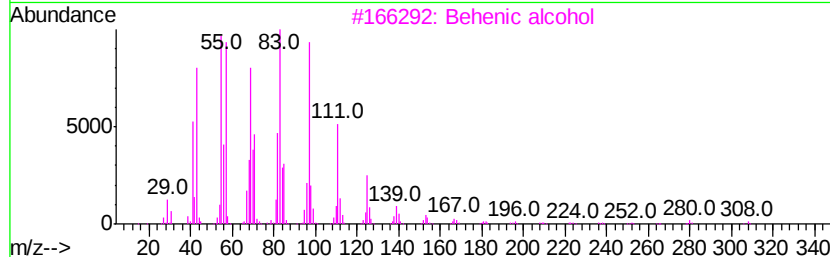
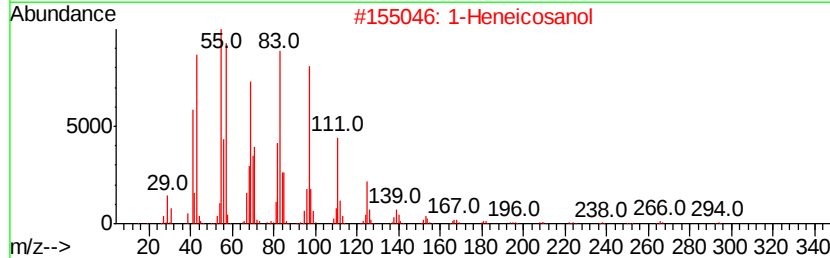
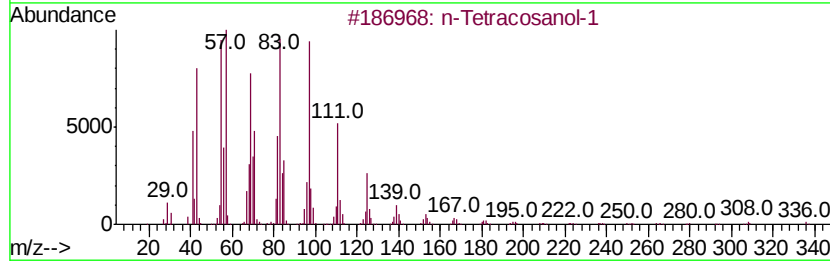
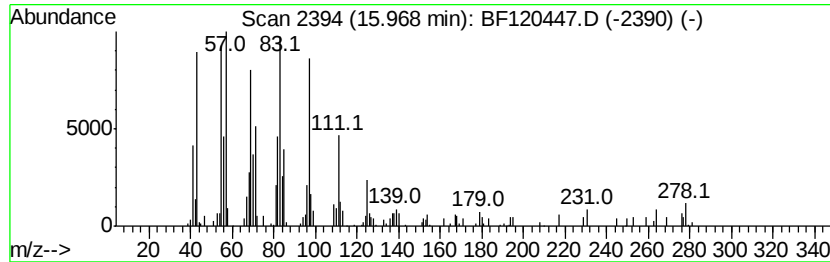
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 n-Tetracosanol-1 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD		R.T.
15.97	3.83 ng	295412	Perylene-d12		15.44
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Tetracosanol-1	354	C24H50O	000506-51-4	95
2	1-Heneicosanol	312	C21H44O	015594-90-8	95
3	Behenic alcohol	326	C22H46O	000661-19-8	95
4	1-Heptacosanol	396	C27H56O	002004-39-9	91
5	Trifluoroacetoxy hexadecane	338	C18H33F3O2	006222-03-3	81



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Misc :
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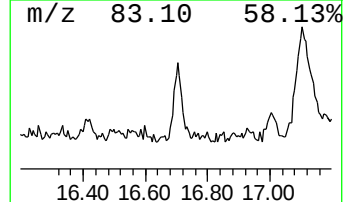
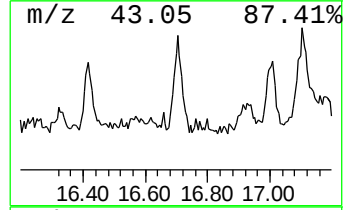
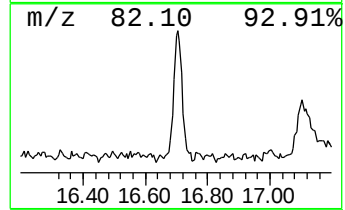
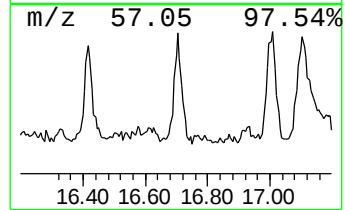
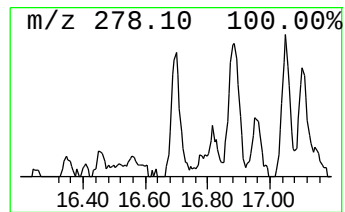
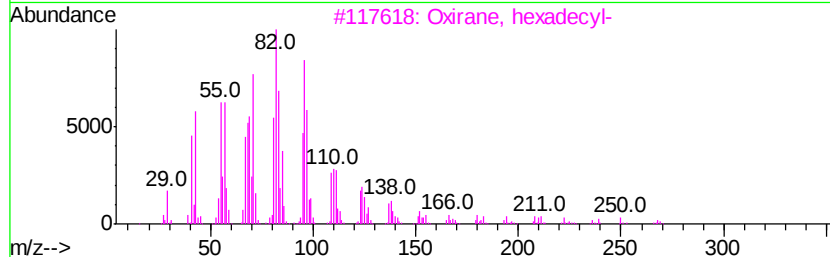
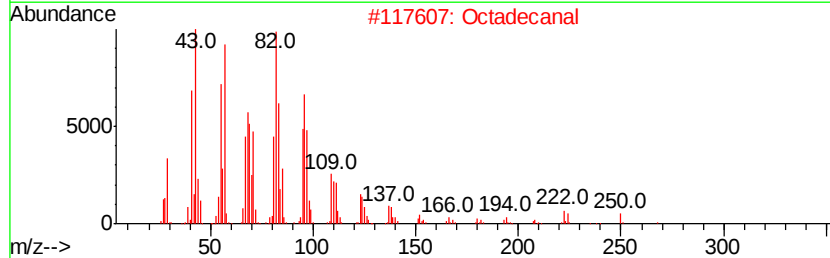
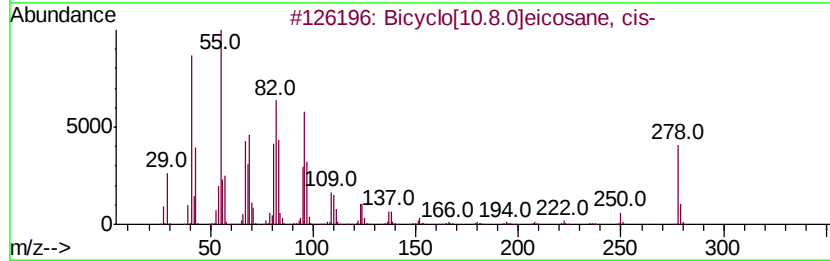
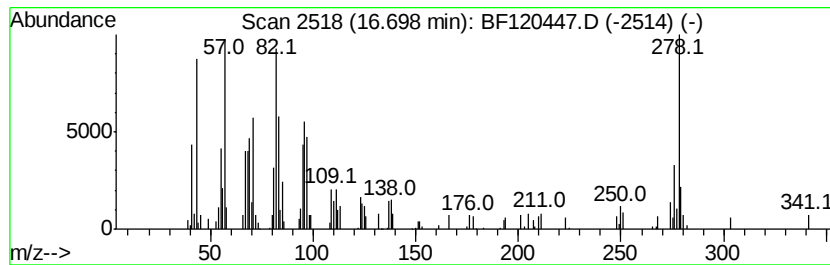
Instrument :
BNA_F
ClientSampleId :
NM59-COMP-2020-0-6

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF052820.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Bicyclo[10.8.0]eicosane, cis- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.	
16.70	2.00 ng	154433	Perylene-d12	15.44	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Bicyclo[10.8.0]eicosane, cis-	278	C20H38	1000155-82-2	93
2	Octadecanal	268	C18H36O	000638-66-4	60
3	Oxirane, hexadecyl-	268	C18H36O	007390-81-0	55
4	1,19-Eicosadiene	278	C20H38	014811-95-1	49
5	11-Dodecen-1-ol trifluoroacetate	280	C14H23F3O2	128792-46-1	46



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF052920\
Data File : BF120447.D
Acq On : 29 May 2020 18:55
Operator : MA/SJ
Sample : L2773-08
Misc :
ALS Vial : 16 Sample Multiplier: 1

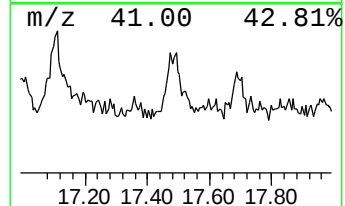
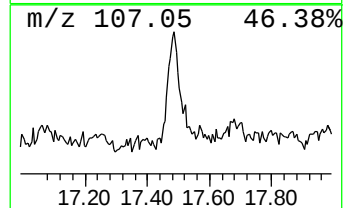
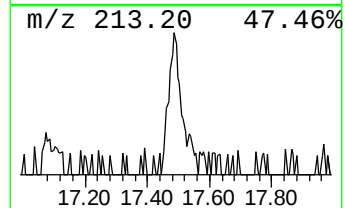
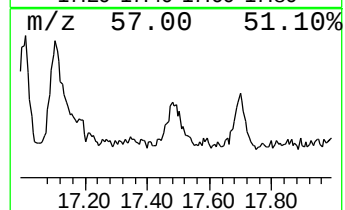
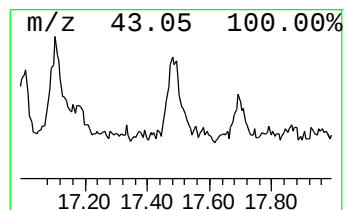
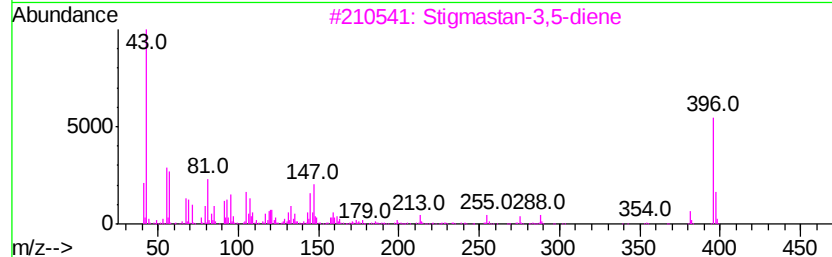
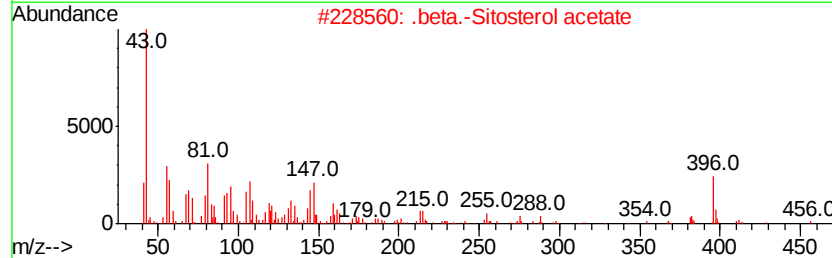
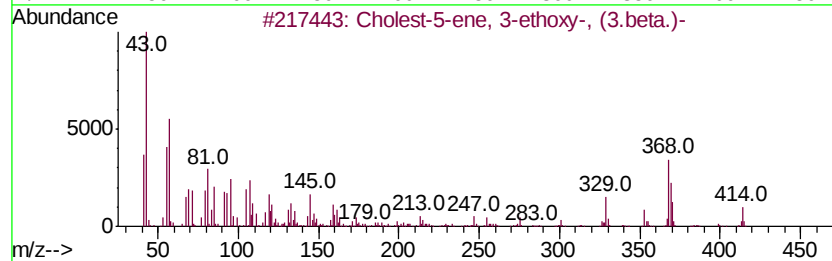
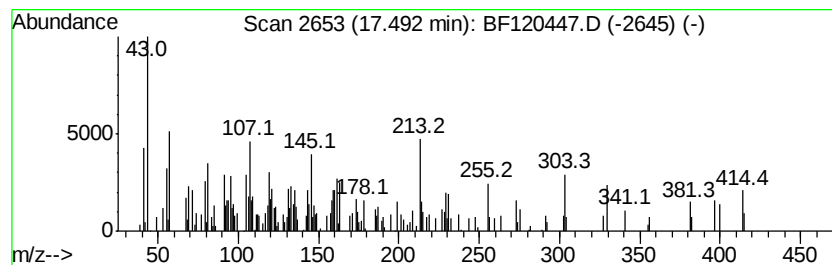
Instrument :
BNA_F
ClientSampleId :
NM59-COMP-2020-0-6

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF052820.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Cholest-5-ene, 3-ethoxy-, (... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD		R.T.		
17.49	4.06 ng	313335	Perylene-d12		15.44		
Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cholest-5-ene, 3-ethoxy-, (3.bet...	414	C29H50O	000986-19-6	62
2			.beta.-Sitosterol acetate	456	C31H52O2	000915-05-9	27
3			Stigmastan-3,5-diene	396	C29H48	1000214-16-4	25
4			Norflurazon	303	C12H9ClF3N3O	027314-13-2	11
5			N-(3-Chloro-4-fluoro-phenyl)-2-(...	329	C14H10Cl2FNOS	1000295-12-7	10



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF052920\
 Data File : BF120447.D
 Acq On : 29 May 2020 18:55
 Operator : MA/SJ
 Sample : L2773-08
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 NM59-COMP-2020-0-6

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF052820.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
Cyclohexane	1.93	6.1 ng		511448	1	6.84	1671580	20.0
Methylene chloride	1.96	25.8 ng		2154210	1	6.84	1671580	20.0
Butane, 2-methoxy...	2.11	17.6 ng		1475070	1	6.84	1671580	20.0
Ethanol, 2-(2-eth...	6.66	7.2 ng		602537	1	6.84	1671580	20.0
2,4,7,9-Tetrameth...	9.31	14.9 ng		1692390	3	9.87	2277000	20.0
1-Heneicosyl formate	13.85	7.6 ng		709371	5	13.99	1856800	20.0
Octacosyl acetate	14.48	3.4 ng		317907	5	13.99	1856800	20.0
Tetracosane	15.11	2.2 ng		166804	6	15.44	1543160	20.0
Benzo[e]pyrene	15.32	2.3 ng		177157	6	15.44	1543160	20.0
Heneicosane	15.92	3.2 ng		249574	6	15.44	1543160	20.0
n-Tetracosanol-1	15.97	3.8 ng		295412	6	15.44	1543160	20.0
Bicyclo[10.8.0]ei...	16.70	2.0 ng		154433	6	15.44	1543160	20.0
Cholest-5-ene, 3-...	17.49	4.1 ng		313335	6	15.44	1543160	20.0