

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF052920\
 Data File : BF120452.D
 Acq On : 29 May 2020 21:12
 Operator : MA/SJ
 Sample : L2792-09
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 NM84-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	3	8	10	rVB	9419636	16790990	100.00%	24.861%
2	1.957	10	12	28	rVB	1912492	2720118	16.20%	4.027%
3	2.087	30	34	51	rVB	1440786	1963674	11.69%	2.907%
4	2.263	61	64	87	rVB	227654	362209	2.16%	0.536%
5	5.451	601	606	609	rBV	3943098	3835546	22.84%	5.679%
6	6.463	772	778	781	rBV	3940576	4242141	25.26%	6.281%
7	6.834	838	841	845	rBV	1833110	1659846	9.89%	2.458%
8	6.992	864	868	871	rBV	4300421	3822710	22.77%	5.660%
9	7.398	931	937	940	rBV	3263470	2943405	17.53%	4.358%
10	8.116	1055	1059	1062	rBV	2333020	2099368	12.50%	3.108%
11	9.198	1237	1243	1246	rBV	5555207	5310606	31.63%	7.863%
12	9.586	1305	1309	1312	rBV	990641	737772	4.39%	1.092%
13	9.875	1353	1358	1361	rBV	2497551	2197491	13.09%	3.254%
14	10.657	1487	1491	1494	rBV	3203772	2754934	16.41%	4.079%
15	11.357	1605	1610	1612	rBV	2277172	1953845	11.64%	2.893%
16	11.380	1612	1614	1617	rVV	450756	350196	2.09%	0.519%
17	11.886	1697	1700	1704	rBV	211498	184816	1.10%	0.274%
18	12.563	1809	1815	1819	rVB2	681047	871197	5.19%	1.290%
19	12.792	1848	1854	1857	rBV	622727	629315	3.75%	0.932%
20	12.945	1875	1880	1883	rVB	4295860	4009147	23.88%	5.936%
21	13.845	2028	2033	2040	rBV3	587004	844713	5.03%	1.251%
22	13.992	2050	2058	2061	rBV2	1894319	2055713	12.24%	3.044%
23	14.015	2061	2062	2065	rVB	297150	189400	1.13%	0.280%
24	14.474	2136	2140	2146	rBV4	194426	310504	1.85%	0.460%
25	15.021	2229	2233	2236	rBV	283364	408518	2.43%	0.605%
26	15.109	2244	2248	2250	rBV2	182972	210480	1.25%	0.312%
27	15.133	2250	2252	2259	rVB2	156805	217417	1.29%	0.322%
28	15.321	2280	2284	2290	rVB	177323	223844	1.33%	0.331%
29	15.380	2290	2294	2300	rBV	237673	289406	1.72%	0.429%
30	15.445	2300	2305	2309	rBV	1257872	1533833	9.13%	2.271%
31	15.921	2381	2386	2390	rBV	253634	370720	2.21%	0.549%
32	15.968	2390	2394	2404	rVV3	188570	364042	2.17%	0.539%
33	16.709	2514	2520	2527	rVB5	125493	250436	1.49%	0.371%
34	16.874	2544	2548	2554	rBV3	90985	180926	1.08%	0.268%

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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	17.098	2580	2586	2600	rvB4	164617	478066	2.85%	0.708%
36	17.480	2647	2651	2660	rvB7	81379	171466	1.02%	0.254%

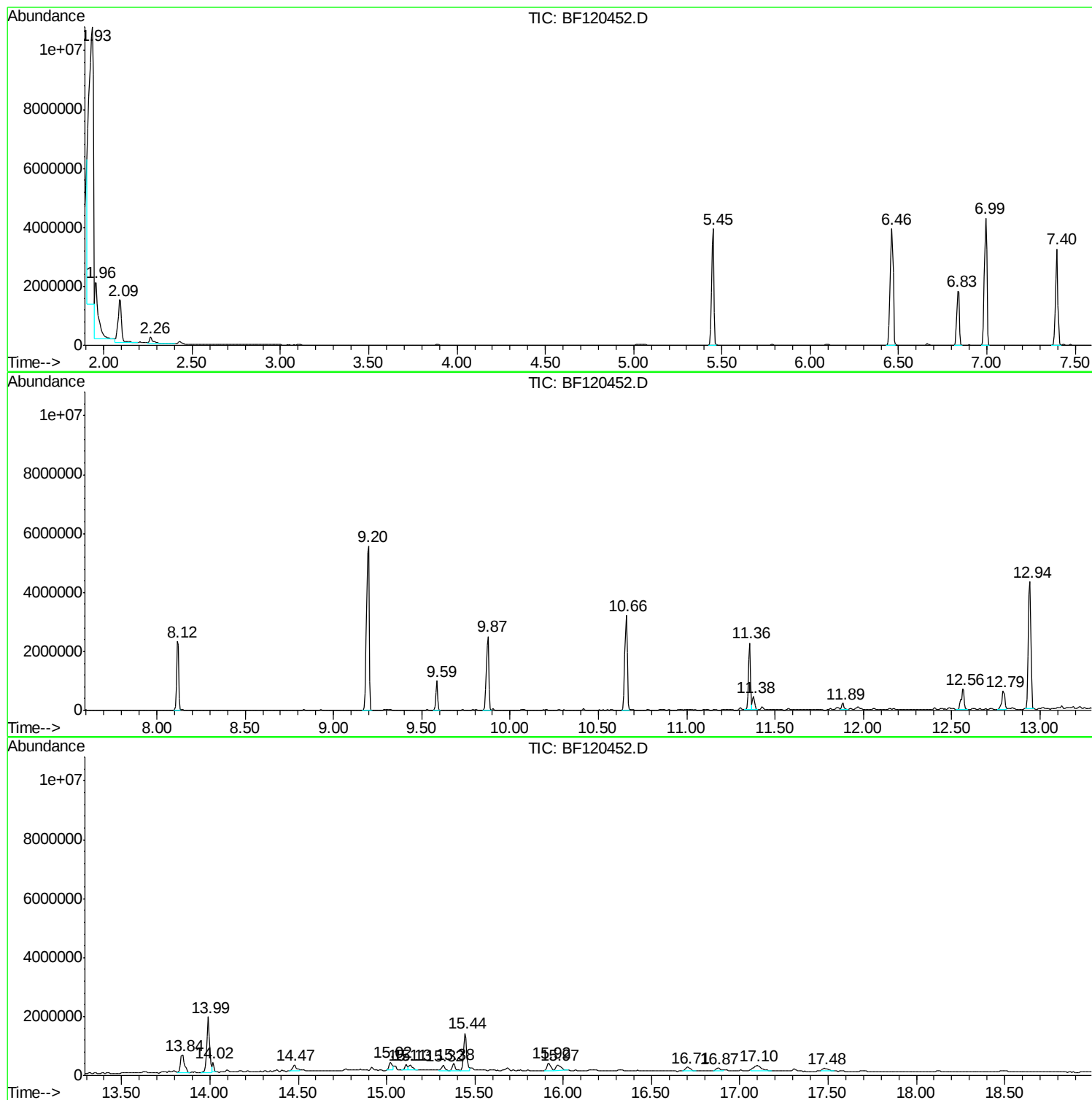
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF052820.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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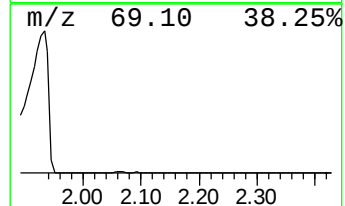
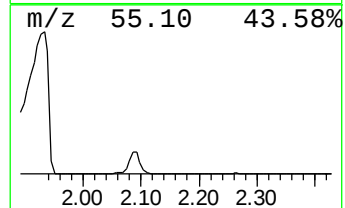
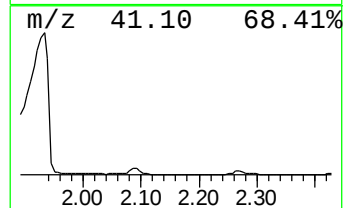
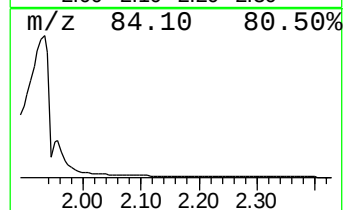
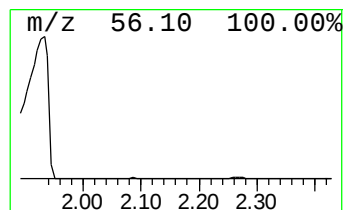
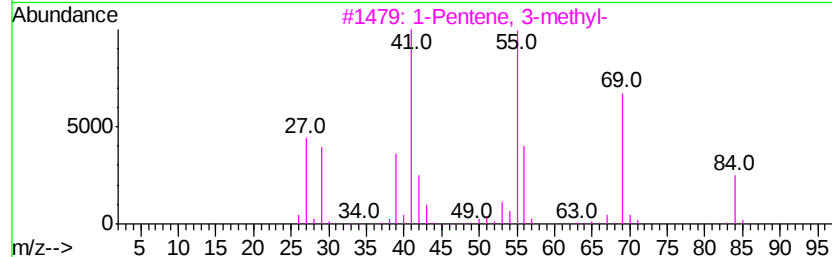
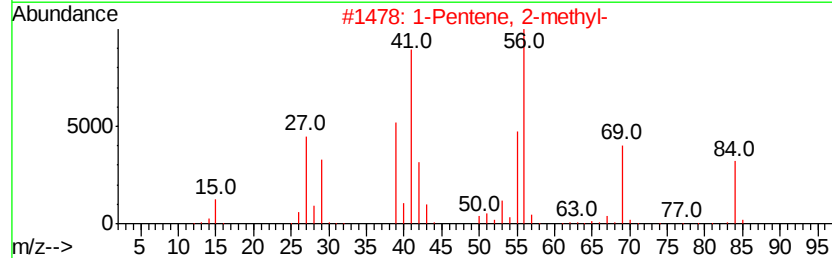
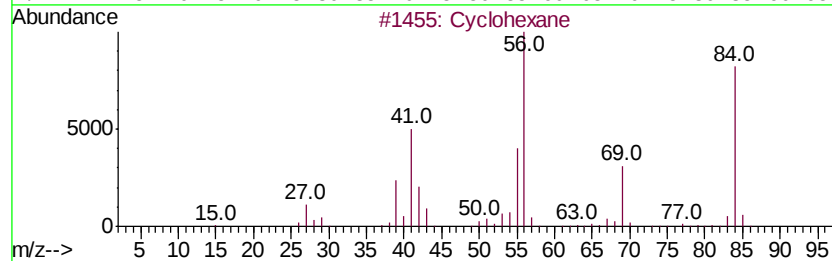
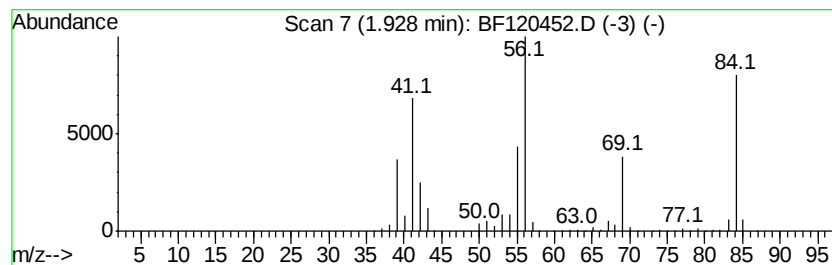
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Cyclohexane Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.93	202.32 ng	16791000	1,4-Dichlorobenzene-d4	6.84
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Cyclohexane	84 C6H12	000110-82-7 93
2		1-Pentene, 2-methyl-	84 C6H12	000763-29-1 86
3		1-Pentene, 3-methyl-	84 C6H12	000760-20-3 43
4		Furan, 2,3-dihydro-4-methyl-	84 C5H8O	034314-83-5 43
5		Pyridine-D5-	84 C5D5N	007291-22-7 42



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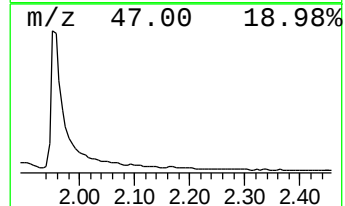
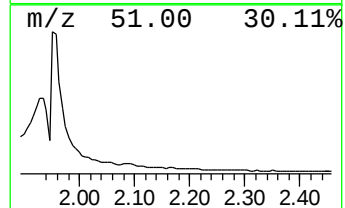
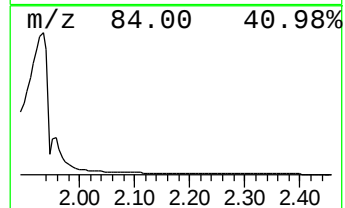
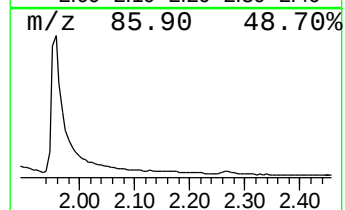
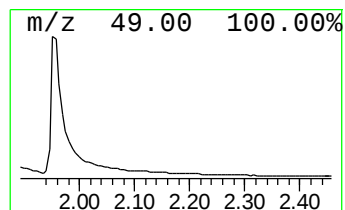
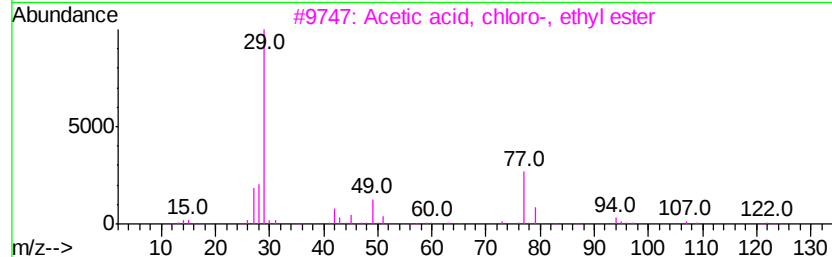
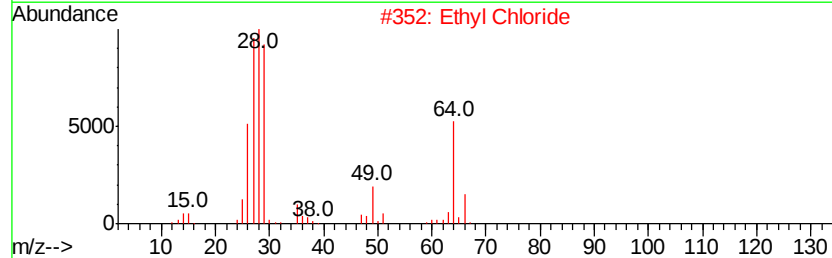
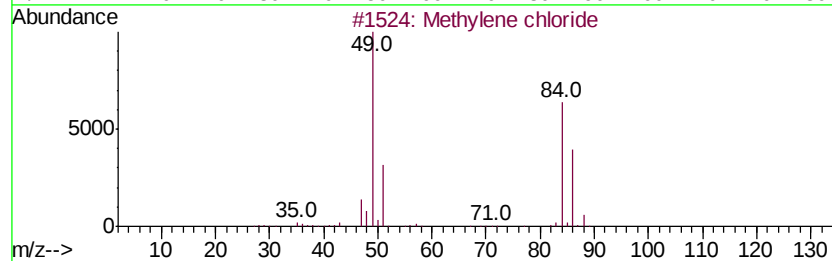
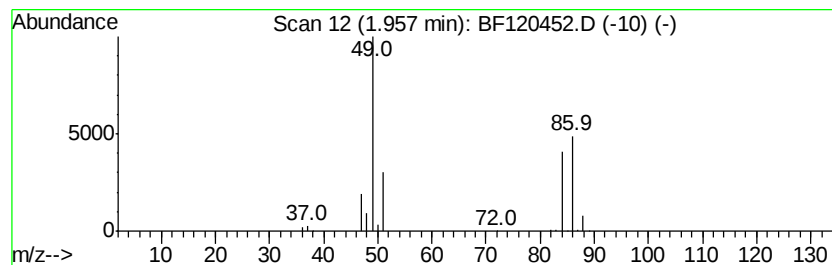
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF052820.M
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Methylene chloride Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.96	32.78 ng	2720120	1,4-Dichlorobenzene-d4	6.84

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Methylene chloride	84	CH2Cl2	000075-09-2	91
2		Ethyl Chloride	64	C2H5Cl	000075-00-3	4
3		Acetic acid, chloro-, ethyl ester	122	C4H7ClO2	000105-39-5	2
4		2-Chloroethanol	80	C2H5ClO	000107-07-3	1
5		Boron trifluoride	68	BF3	007637-07-2	1



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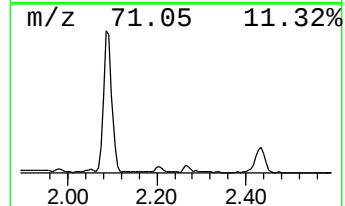
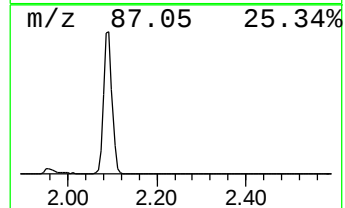
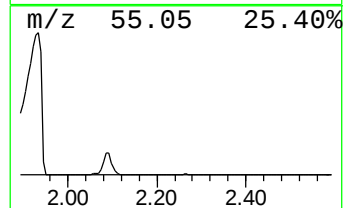
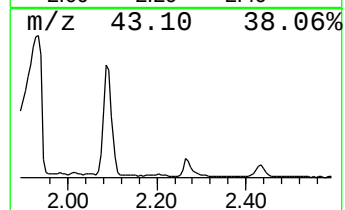
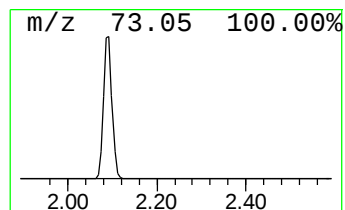
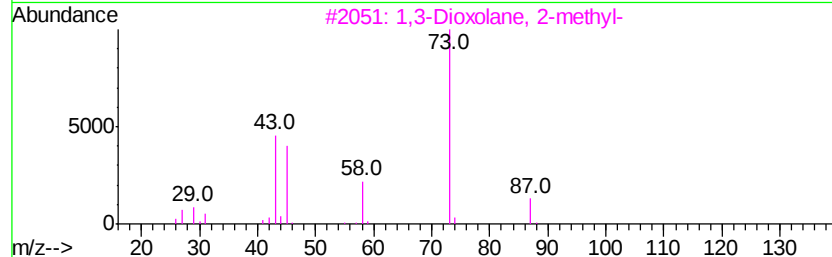
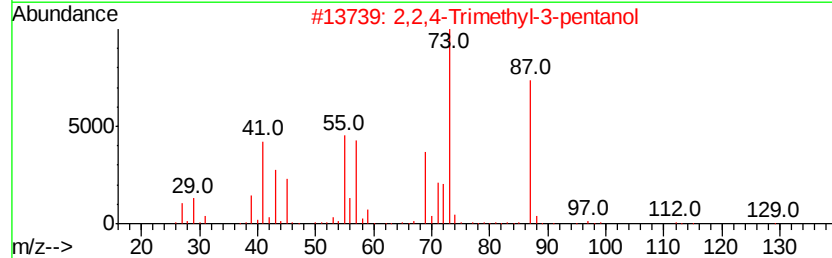
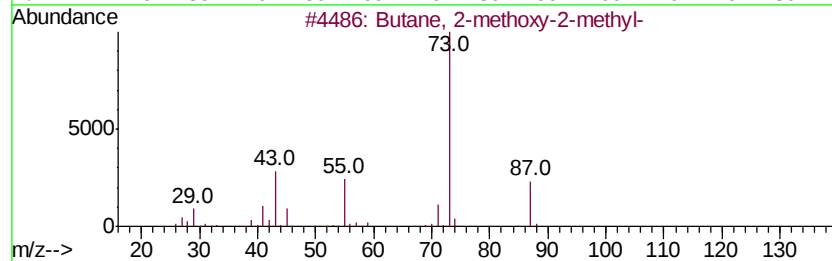
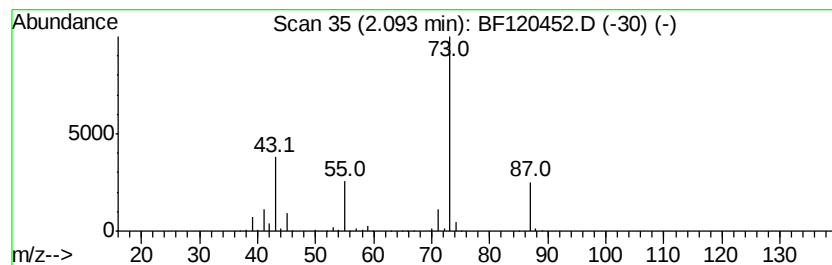
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Butane, 2-methoxy-2-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.	
2.09	23.66 ng	1963670	1,4-Dichlorobenzene-d4	6.84	
Hit# of	5	Tentative ID	MW MolForm	CAS#	Qual
1	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2	2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39
3	1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	25
4	Pentane, 3-methoxy-	102	C6H14O	036839-67-5	12
5	Silane, tetramethyl-	88	C4H12Si	000075-76-3	10



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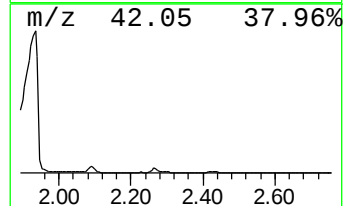
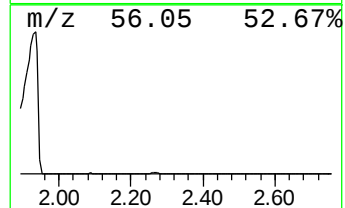
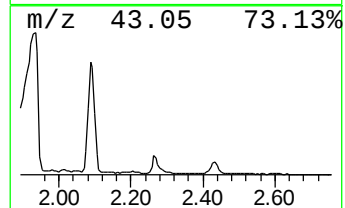
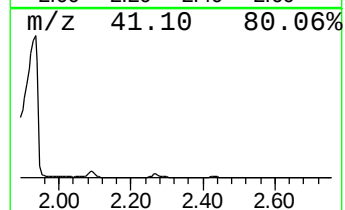
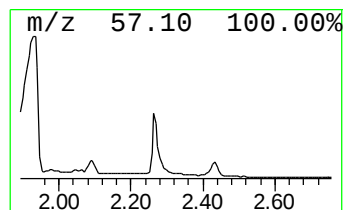
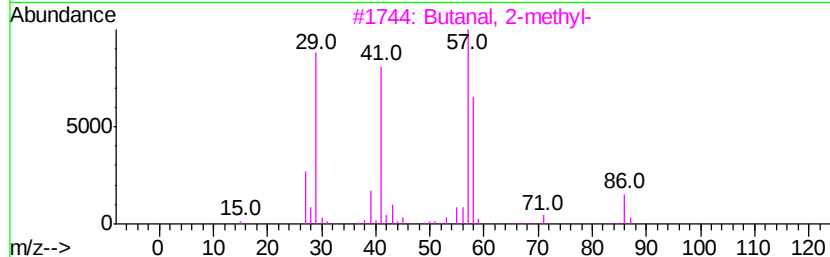
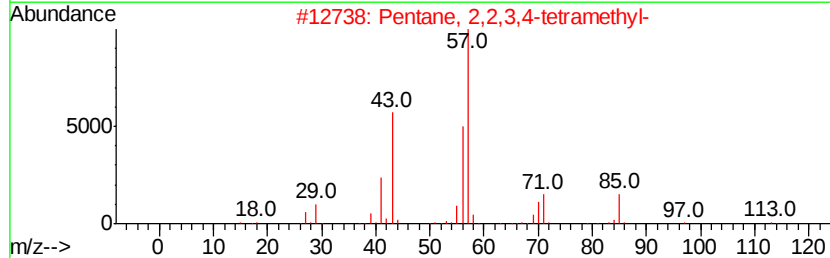
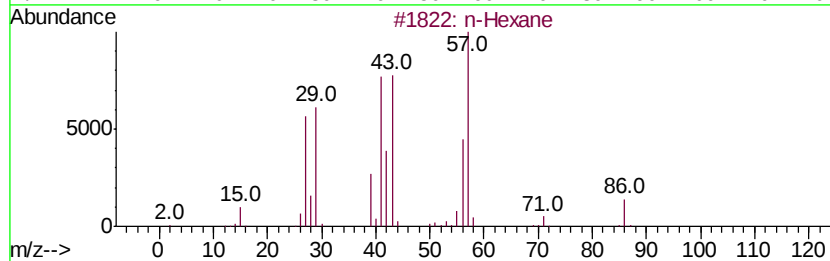
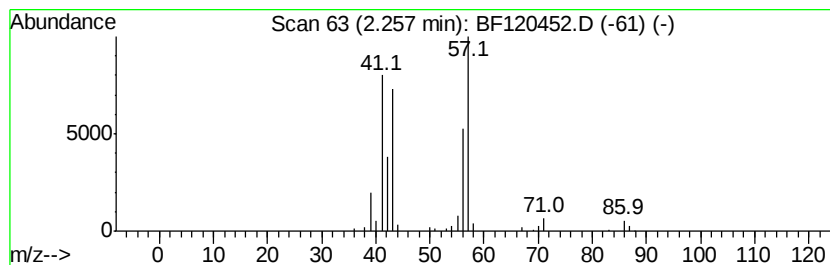
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 n-Hexane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.26	4.36 ng	362209	1,4-Dichlorobenzene-d4	6.84

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexane	86	C6H14	000110-54-3	91
2			Pentane, 2,2,3,4-tetramethyl-	128	C9H20	001186-53-4	38
3			Butanal, 2-methyl-	86	C5H10O	000096-17-3	38
4			1-Pentene, 4-methyl-	84	C6H12	000691-37-2	33
5			3-Buten-1-ol, 3-methyl-	86	C5H10O	000763-32-6	28



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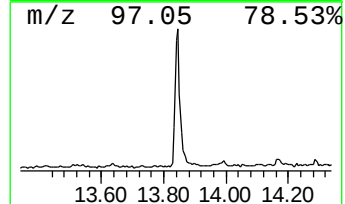
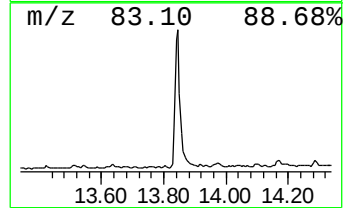
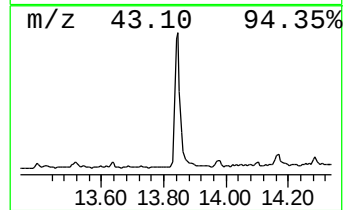
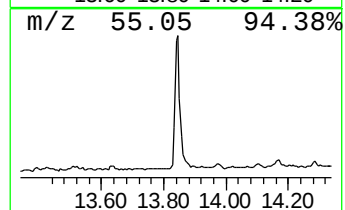
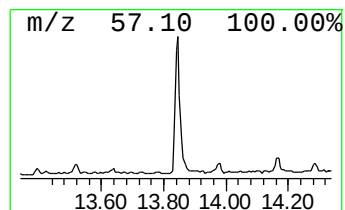
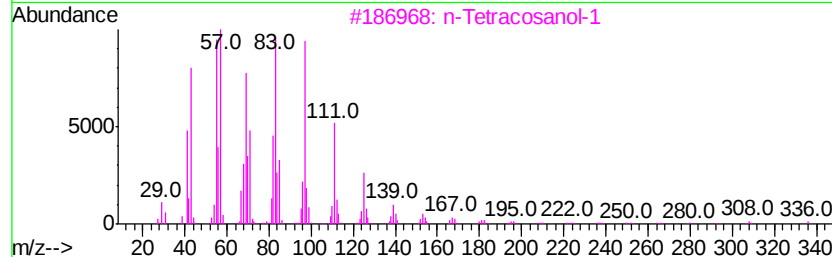
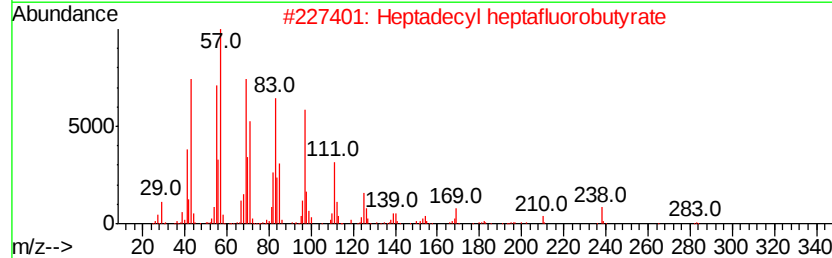
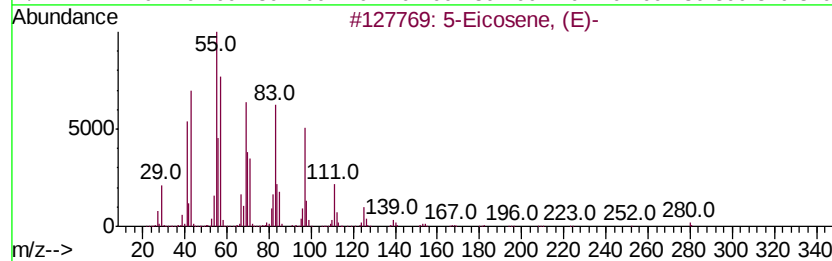
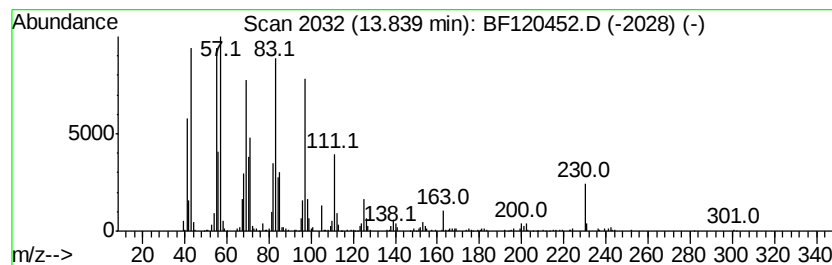
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 5-Eicosene, (E)- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.84	8.22 ng	844713	Chrysene-d12	13.99	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	5-Eicosene, (E)-	280	C20H40	074685-30-6	97
2	Heptadecyl heptafluorobutyrte	452	C21H35F7O2	959085-66-6	94
3	n-Tetracosanol-1	354	C24H50O	000506-51-4	94
4	1-Heneicosanol	312	C21H44O	015594-90-8	93
5	Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF052920\
Data File : BF120452.D
Acq On : 29 May 2020 21:12
Operator : MA/SJ
Sample : L2792-09
Misc :
ALS Vial : 21 Sample Multiplier: 1

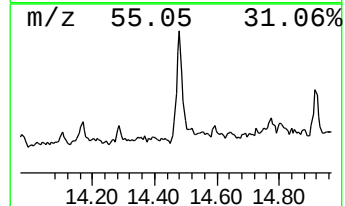
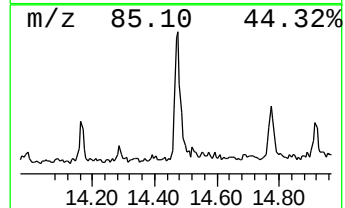
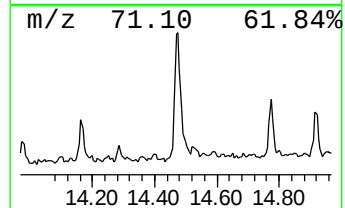
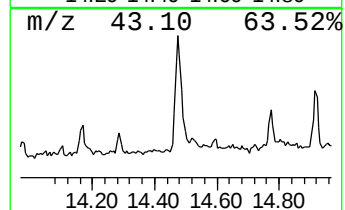
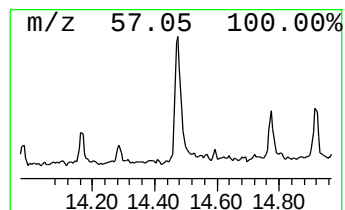
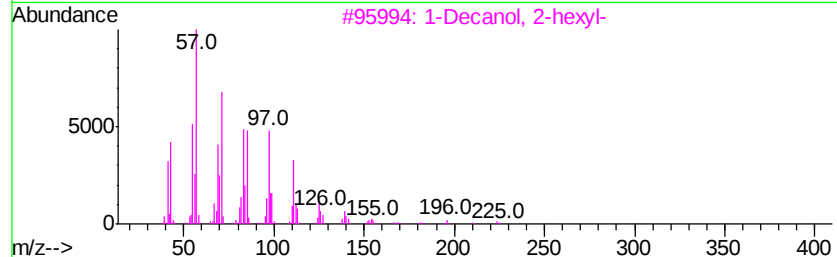
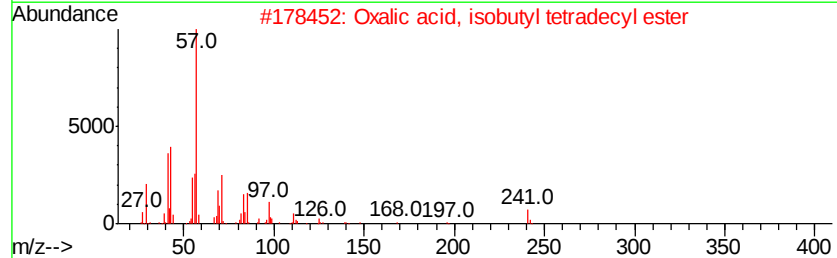
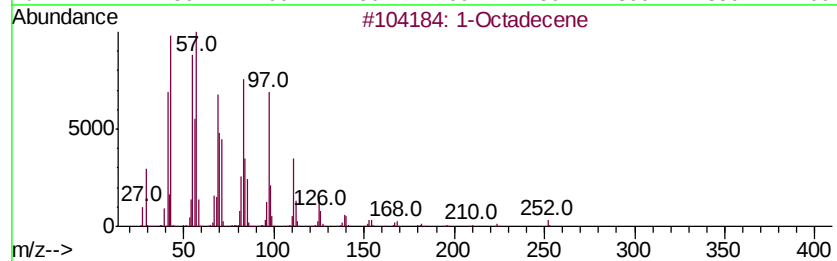
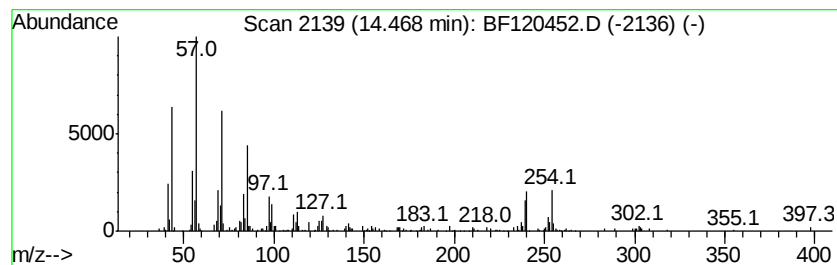
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ClientSampleId :
NM84-COMP-2020-0-6

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 7 1-Octadecene Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.47	3.02 ng	310504	Chrysene-d12	13.99
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	1-Octadecene		252 C18H36	000112-88-9 95
2	Oxalic acid, isobutyl tetradecyl...		342 C20H38O4	1000309-37-9 81
3	1-Decanol, 2-hexyl-		242 C16H34O	002425-77-6 81
4	Oxalic acid, isobutyl hexadecyl ...		370 C22H42O4	1000309-38-1 74
5	1-Hexadecanol, 2-methyl-		256 C17H36O	002490-48-4 74



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF052920\
 Data File : BF120452.D
 Acq On : 29 May 2020 21:12
 Operator : MA/SJ
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Instrument :
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 ClientSampleId :
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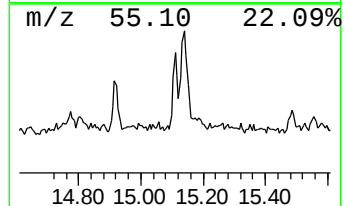
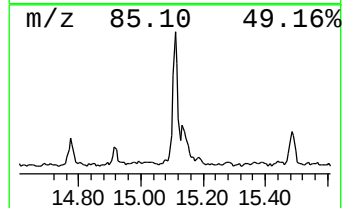
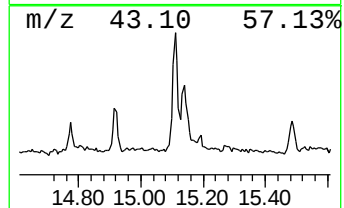
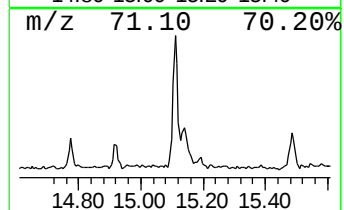
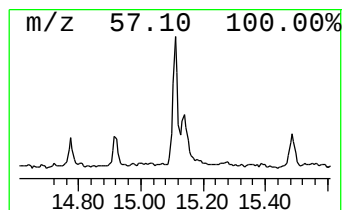
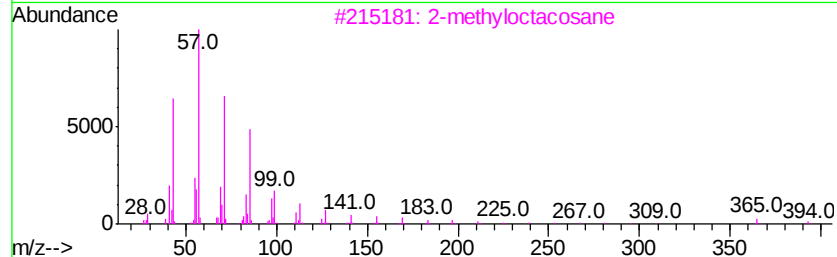
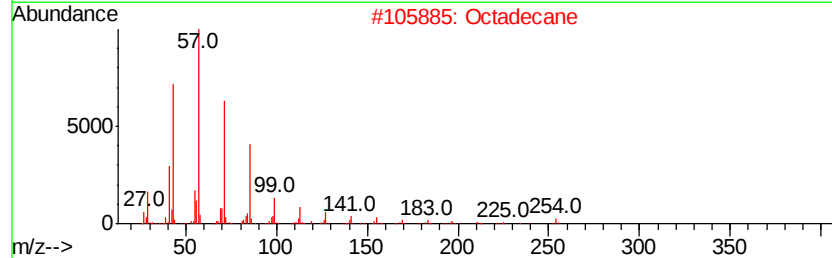
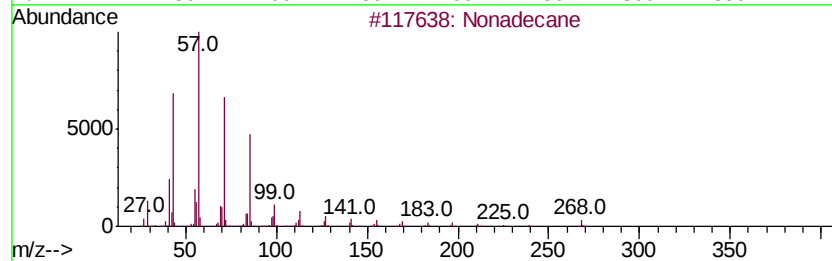
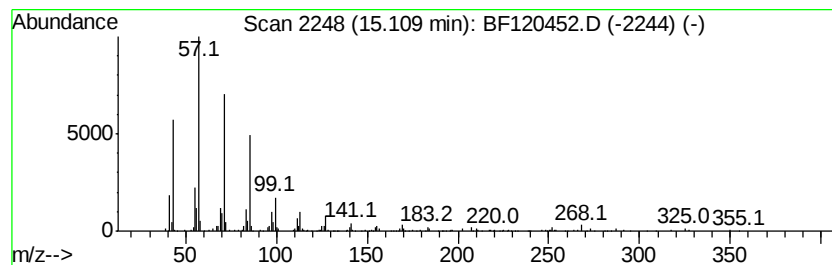
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

 Peak Number 8 Nonadecane Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.11	2.74 ng	210480	Perylene-d12	15.44

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonadecane		268	C19H40	000629-92-5	94
2	Octadecane		254	C18H38	000593-45-3	91
3	2-methyloctacosane		408	C29H60	1000376-72-8	87
4	Heneicosane		296	C21H44	000629-94-7	87
5	Octacosane		394	C28H58	000630-02-4	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF052920\
Data File : BF120452.D
Acq On : 29 May 2020 21:12
Operator : MA/SJ
Sample : L2792-09
Misc :
ALS Vial : 21 Sample Multiplier: 1

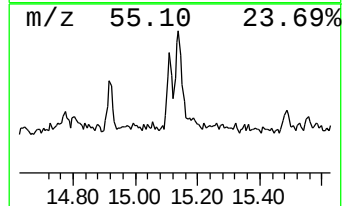
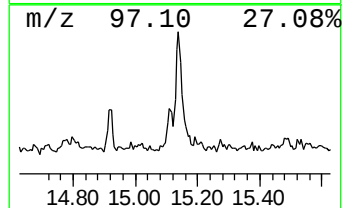
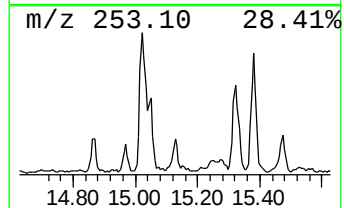
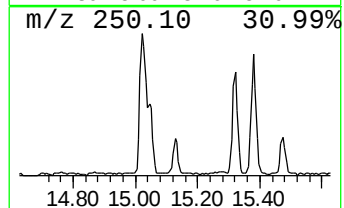
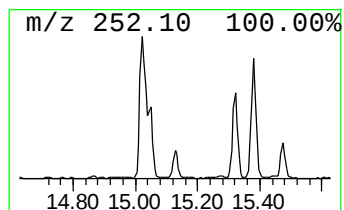
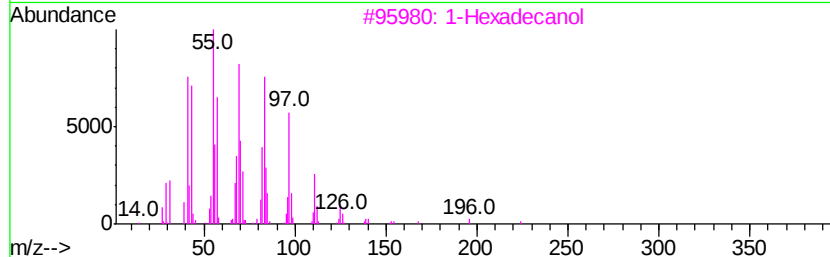
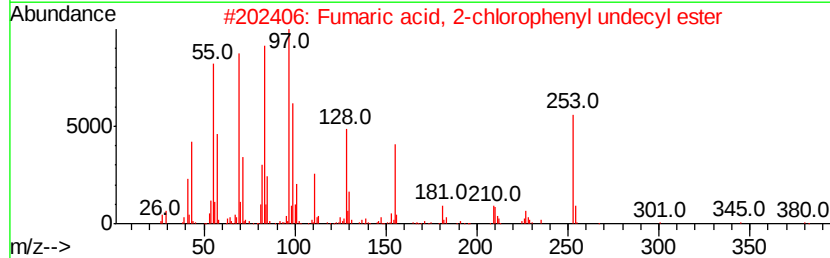
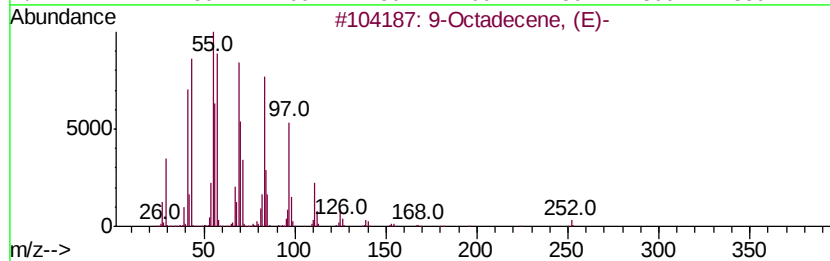
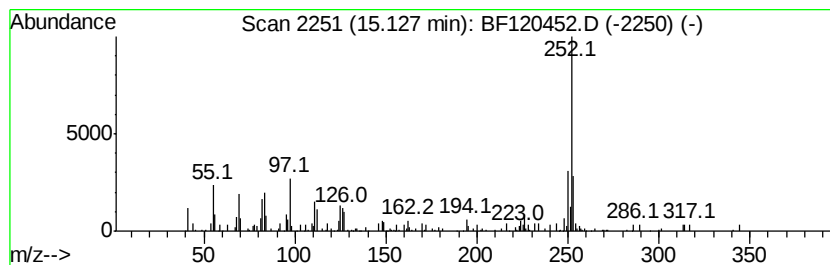
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF052820.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 9-Octadecene, (E)- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.13	2.83 ng	217417	Perylene-d12	15.44
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	9-Octadecene, (E)-	252 C18H36	007206-25-9	58
2	Fumaric acid, 2-chlorophenyl und...	380 C21H29ClO4	1000344-83-3	46
3	1-Hexadecanol	242 C16H34O	036653-82-4	25
4	Benzo[a]pyrene	252 C20H12	000050-32-8	25
5	2,5-Dimethyl-1-pyrroline	97 C6H11N	000822-64-0	25



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF052920\
Data File : BF120452.D
Acq On : 29 May 2020 21:12
Operator : MA/SJ
Sample : L2792-09
Misc :
ALS Vial : 21 Sample Multiplier: 1

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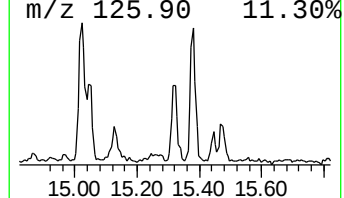
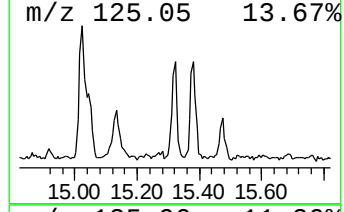
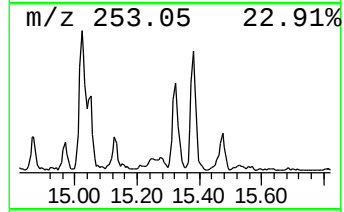
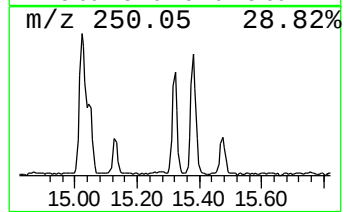
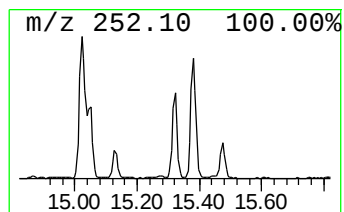
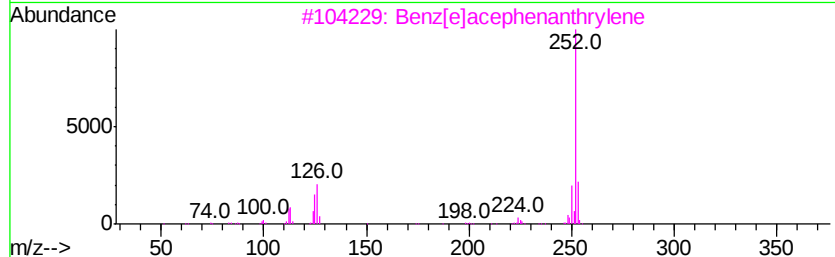
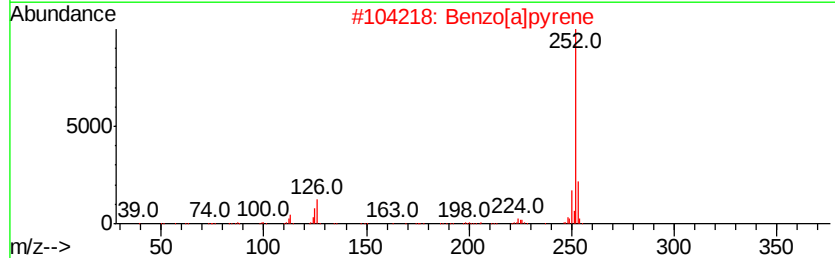
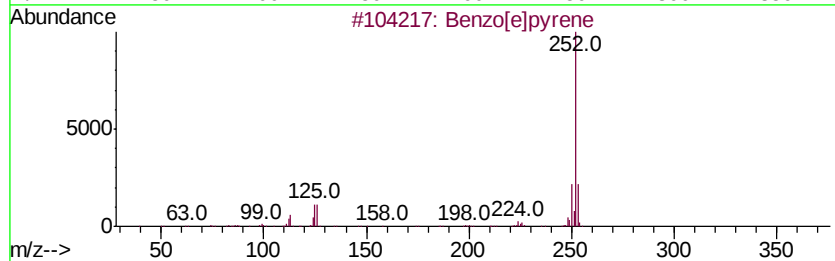
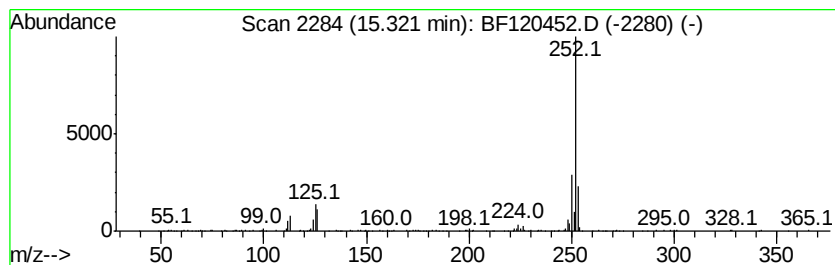
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Benzo[e]pyrene Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.32	2.92 ng	223844	Perylene-d12	15.44

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzo[e]pyrene	252	C20H12	000192-97-2	98
2		Benzo[a]pyrene	252	C20H12	000050-32-8	96
3		Benz[e]acephenanthrylene	252	C20H12	000205-99-2	95
4		Perylene	252	C20H12	000198-55-0	90
5		Benzo[k]fluoranthene	252	C20H12	000207-08-9	89



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF052920\
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 Acq On : 29 May 2020 21:12
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Instrument :
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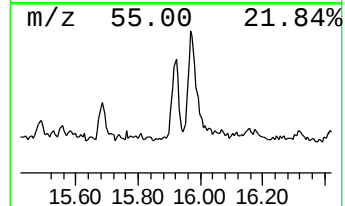
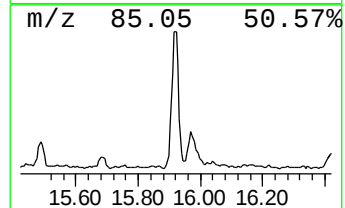
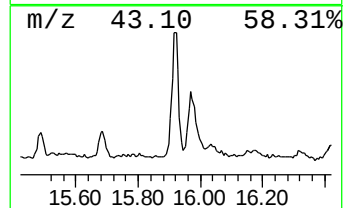
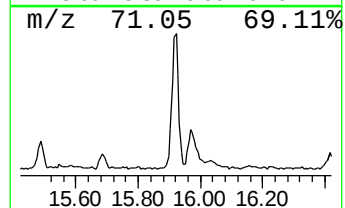
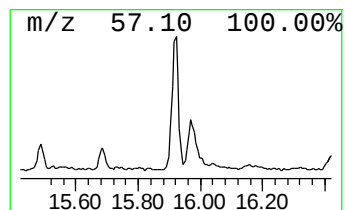
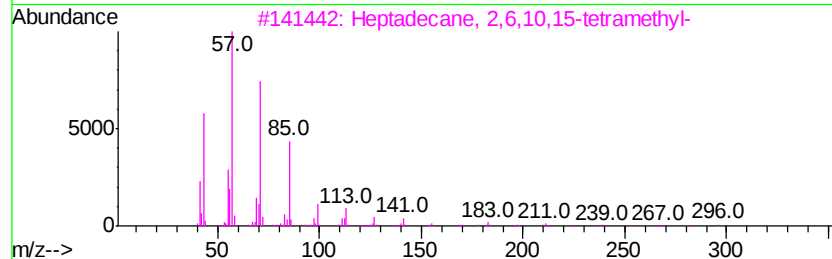
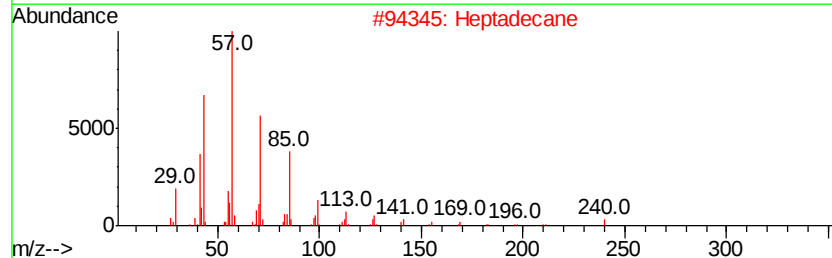
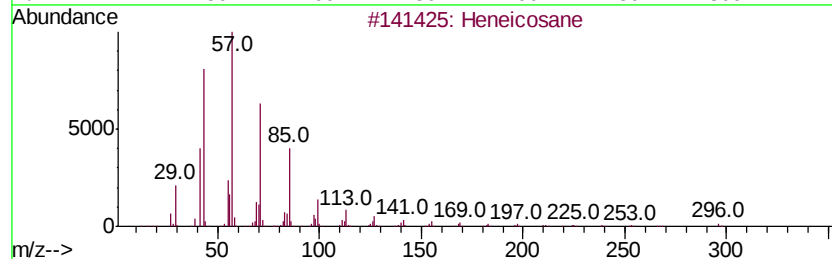
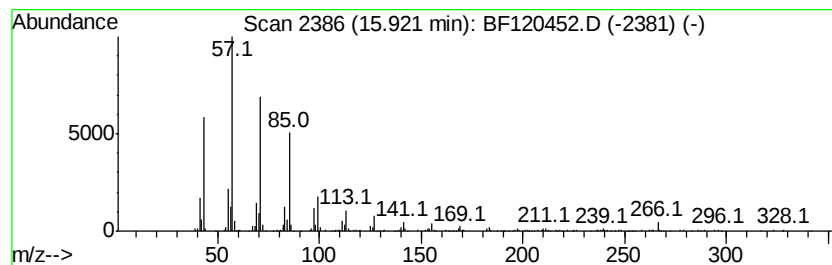
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TIC Library : C:\DATABASE\NIST11.L
 TIC Integration Parameters: LSCINT.P

 Peak Number 11 Heneicosane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.92	4.83 ng	370720	Perylene-d12	15.44

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosane	296	C21H44	000629-94-7	96
2		Heptadecane	240	C17H36	000629-78-7	93
3		Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	93
4		Docosane, 11-butyl-	366	C26H54	013475-76-8	91
5		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91



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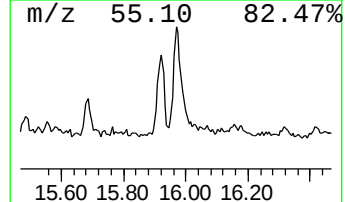
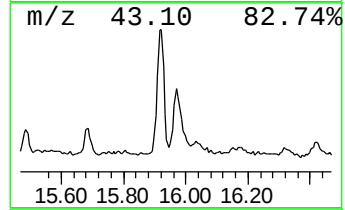
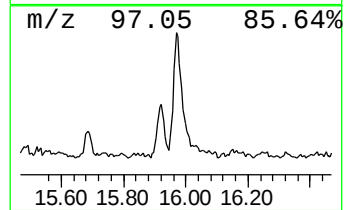
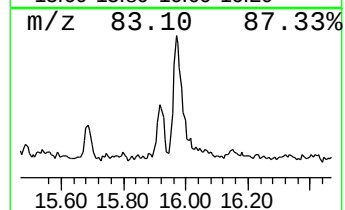
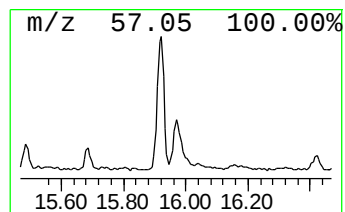
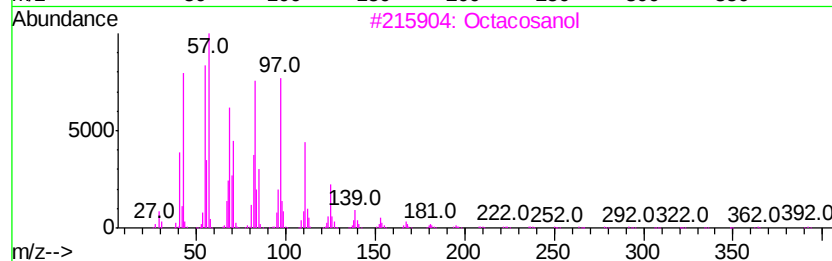
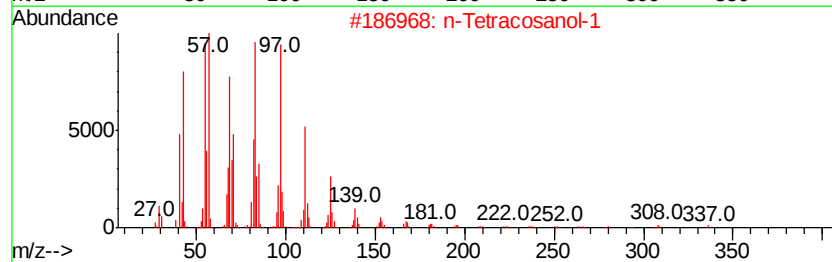
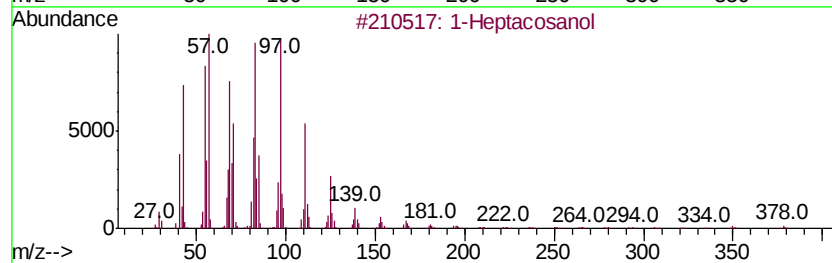
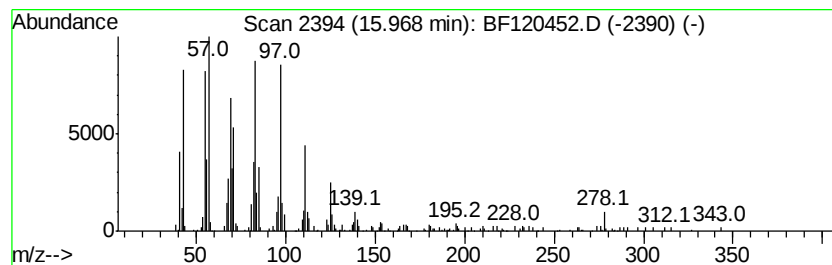
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 1-Heptacosanol Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD		R.T.
15.97	4.75 ng	364042	Perylene-d12		15.44
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Heptacosanol	396	C27H56O	002004-39-9	95
2	n-Tetracosanol-1	354	C24H50O	000506-51-4	95
3	Octacosanol	410	C28H58O	000557-61-9	95
4	Octacosyl acetate	452	C30H60O2	018206-97-8	94
5	Heptacosyl acetate	438	C29H58O2	1000351-78-2	93



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF052920\
Data File : BF120452.D
Acq On : 29 May 2020 21:12
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ClientSampleId :
NM84-COMP-2020-0-6

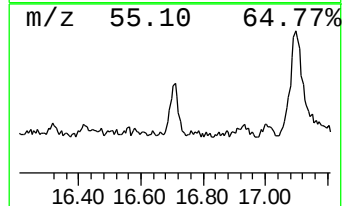
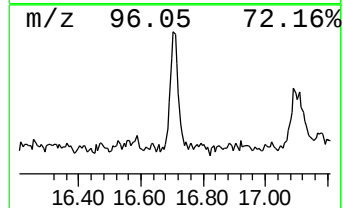
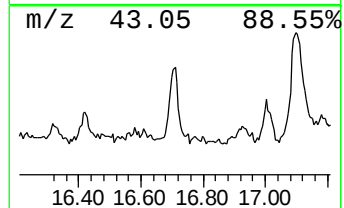
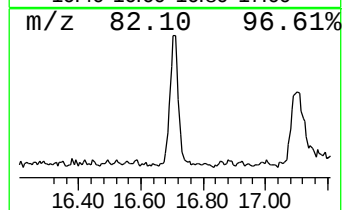
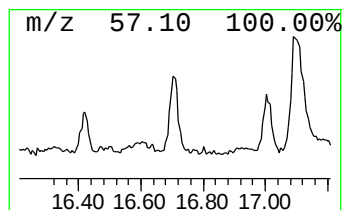
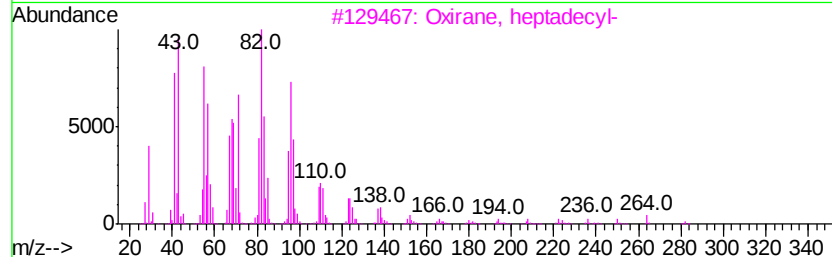
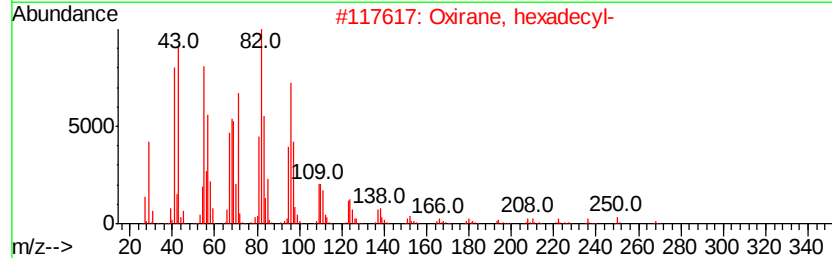
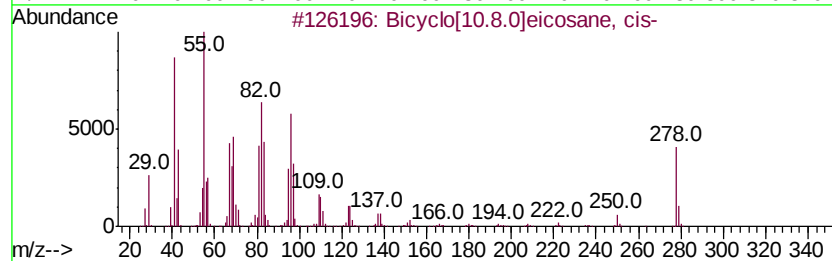
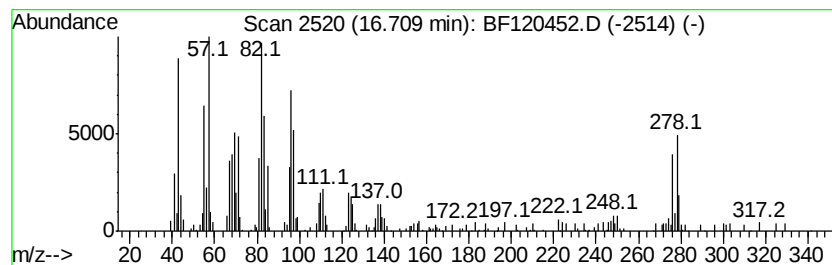
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 13 Bicyclo[10.8.0]eicosane, cis- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.71	3.27 ng	250436	Perylene-d12	15.44

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Bicyclo[10.8.0]eicosane, cis-	278	C20H38	1000155-82-2	64
2		Oxirane, hexadecyl-	268	C18H36O	007390-81-0	49
3		Oxirane, heptadecyl-	282	C19H38O	067860-04-2	49
4		Cyclohexane, 1-(1,5-dimethylhexyl)-	280	C20H40	056009-20-2	46
5		1,22-Docosanediol	342	C22H46O2	022513-81-1	43



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF052920\
Data File : BF120452.D
Acq On : 29 May 2020 21:12
Operator : MA/SJ
Sample : L2792-09
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
NM84-COMP-2020-0-6

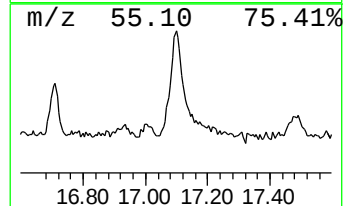
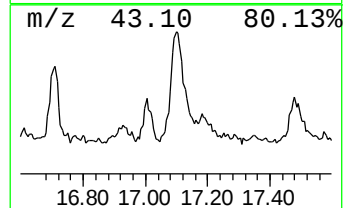
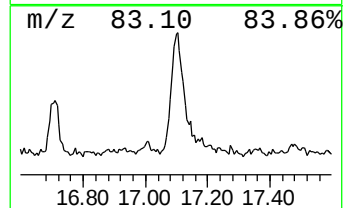
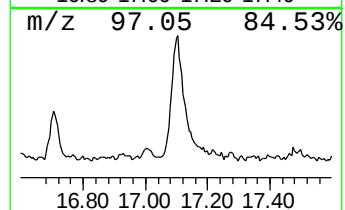
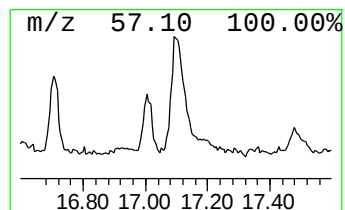
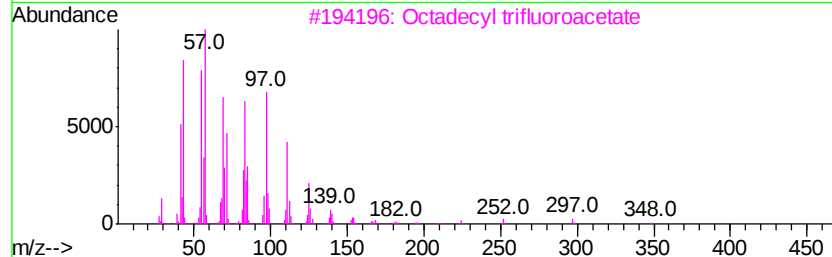
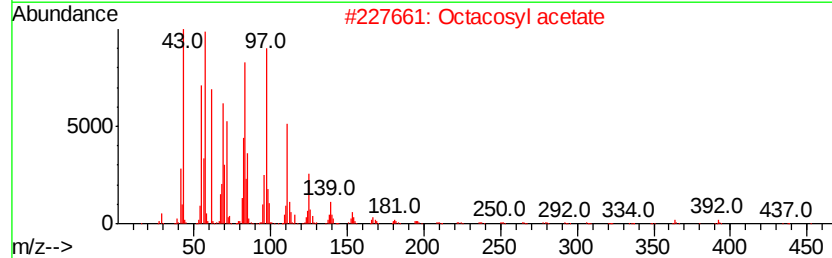
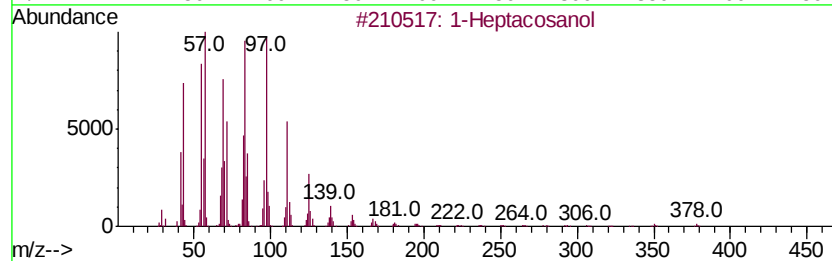
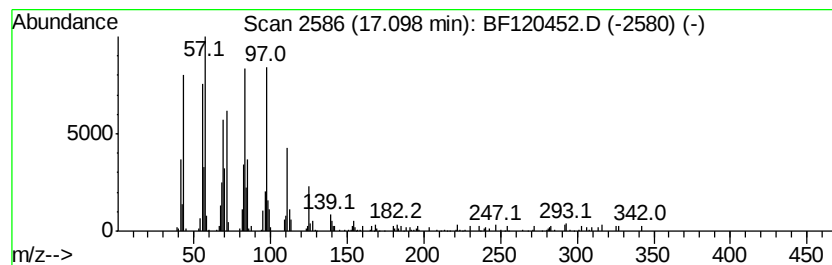
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Octacosyl acetate Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.10	6.23 ng	478066	Perylene-d12	15.44

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		1-Heptacosanol	396	C27H56O	002004-39-9	95
2		Octacosyl acetate	452	C30H60O2	018206-97-8	95
3		Octadecyl trifluoroacetate	366	C20H37F3O2	079392-43-1	94
4		Pentafluoropropionic acid, hepta...	402	C20H35F5O2	959218-78-1	94
5		Pentafluoropropionic acid, octad...	416	C21H37F5O2	959261-25-7	94



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF052920\
Data File : BF120452.D
Acq On : 29 May 2020 21:12
Operator : MA/SJ
Sample : L2792-09
Misc :
ALS Vial : 21 Sample Multiplier: 1

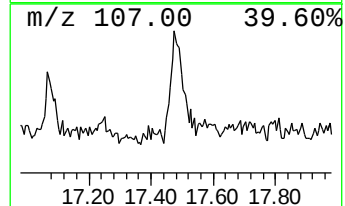
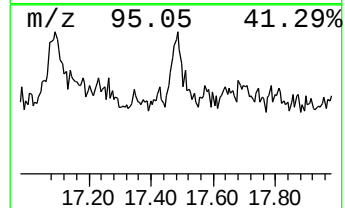
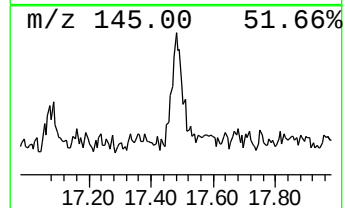
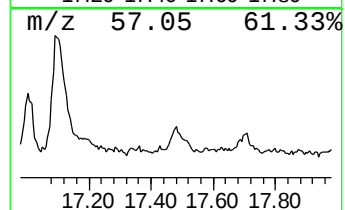
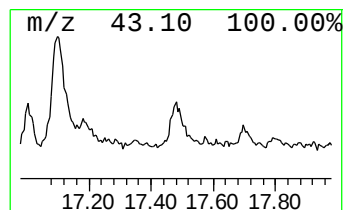
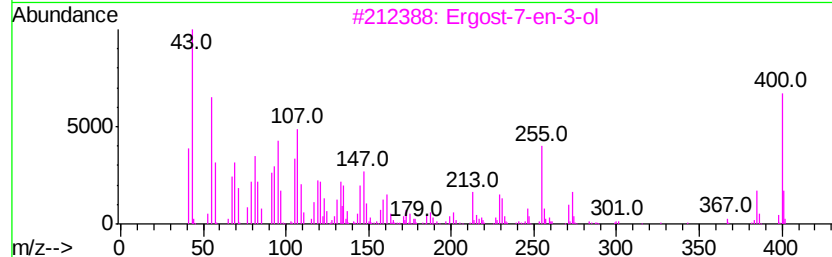
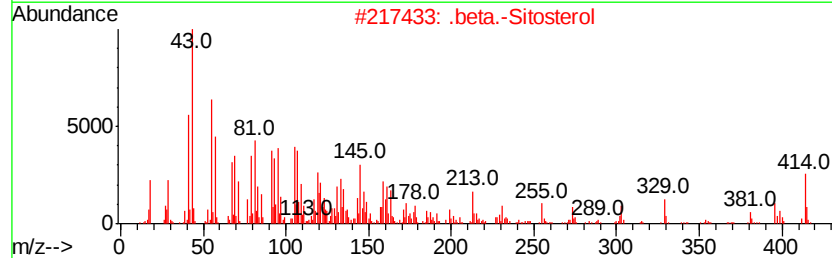
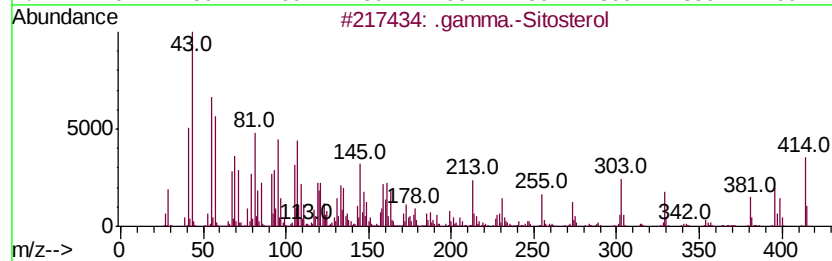
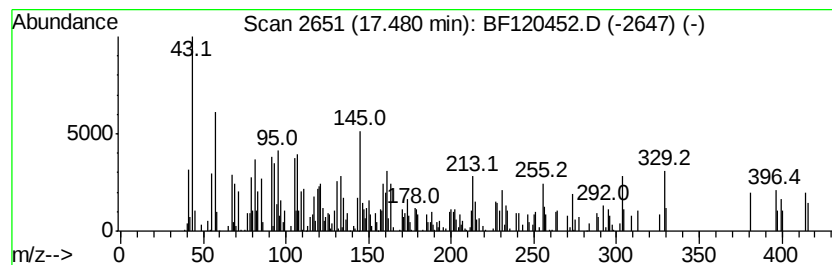
Instrument :
BNA_F
ClientSampleId :
NM84-COMP-2020-0-6

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 15 .gamma.-Sitosterol Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.	
17.48	2.24 ng	171466	Perylene-d12	15.44	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	.gamma.-Sitosterol	414	C29H50O	000083-47-6	70
2	.beta.-Sitosterol	414	C29H50O	000083-46-5	60
3	Ergost-7-en-3-ol	400	C28H48O	116179-22-7	51
4	Cholest-5-ene, 3-ethoxy-, (3.bet...	414	C29H50O	000986-19-6	46
5	Stigmastan-3-en-6-ol	414	C29H50O	1000214-19-7	41



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF052920\
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 Acq On : 29 May 2020 21:12
 Operator : MA/SJ
 Sample : L2792-09
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 NM84-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	202.3	ng	16791000	1	6.84	1659850	20.0
Methylene chloride	1.96	32.8	ng	2720120	1	6.84	1659850	20.0
Butane, 2-methoxy...	2.09	23.7	ng	1963670	1	6.84	1659850	20.0
n-Hexane	2.26	4.4	ng	362209	1	6.84	1659850	20.0
5-Eicosene, (E)-	13.84	8.2	ng	844713	5	13.99	2055710	20.0
1-Octadecene	14.47	3.0	ng	310504	5	13.99	2055710	20.0
Nonadecane	15.11	2.7	ng	210480	6	15.44	1533830	20.0
9-Octadecene, (E)-	15.13	2.8	ng	217417	6	15.44	1533830	20.0
Benzo[e]pyrene	15.32	2.9	ng	223844	6	15.44	1533830	20.0
Heneicosane	15.92	4.8	ng	370720	6	15.44	1533830	20.0
1-Heptacosanol	15.97	4.8	ng	364042	6	15.44	1533830	20.0
Bicyclo[10.8.0]ei...	16.71	3.3	ng	250436	6	15.44	1533830	20.0
Octacosyl acetate	17.10	6.2	ng	478066	6	15.44	1533830	20.0
.gamma.-Sitosterol	17.48	2.2	ng	171466	6	15.44	1533830	20.0