

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF121719\
 Data File : BF118239.D
 Acq On : 17 Dec 2019 21:16
 Operator : JU
 Sample : K6324-05
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 EP-13

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-BF120619.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.128	3	7	18	rVB	229608	310234	10.84%	1.262%
2	5.069	503	507	518	rVB	326245	411989	14.39%	1.676%
3	5.481	572	577	580	rBV	2475842	2265214	79.12%	9.214%
4	6.492	744	749	752	rBV	2360903	2415865	84.38%	9.827%
5	6.634	768	773	776	rBV	2867670	2646843	92.45%	10.767%
6	6.857	808	811	817	rBV	814408	691219	24.14%	2.812%
7	7.016	834	838	841	rBV	2513190	2077169	72.55%	8.450%
8	7.428	903	908	922	rBV	1414128	1520978	53.12%	6.187%
9	8.145	1026	1030	1051	rBV	935492	865404	30.23%	3.520%
10	9.228	1209	1214	1217	rBV	3150174	2863087	100.00%	11.647%
11	9.622	1277	1281	1289	rVB	109354	102297	3.57%	0.416%
12	9.910	1326	1330	1342	rBV	1043201	919052	32.10%	3.739%
13	10.704	1461	1465	1481	rBV	1594417	1604471	56.04%	6.527%
14	11.404	1580	1584	1593	rBV	846767	797330	27.85%	3.243%
15	11.922	1669	1672	1681	rBV	144531	168333	5.88%	0.685%
16	12.716	1804	1807	1814	rBV3	25172	34804	1.22%	0.142%
17	12.992	1850	1854	1857	rBV	2815273	2399148	83.80%	9.759%
18	13.886	2002	2006	2017	rBV	235322	381694	13.33%	1.553%
19	14.057	2031	2035	2046	rVB	749323	769120	26.86%	3.129%
20	14.210	2058	2061	2064	rBV	35942	39332	1.37%	0.160%
21	14.515	2109	2113	2115	rBV	69301	81724	2.85%	0.332%
22	14.821	2163	2165	2169	rVB	71320	66504	2.32%	0.271%
23	15.168	2220	2224	2228	rBV	100123	115054	4.02%	0.468%
24	15.545	2283	2288	2299	rBV	511095	793377	27.71%	3.227%
25	16.004	2362	2366	2371	rVB	99281	144955	5.06%	0.590%
26	16.521	2450	2454	2458	rVB3	64745	98012	3.42%	0.399%

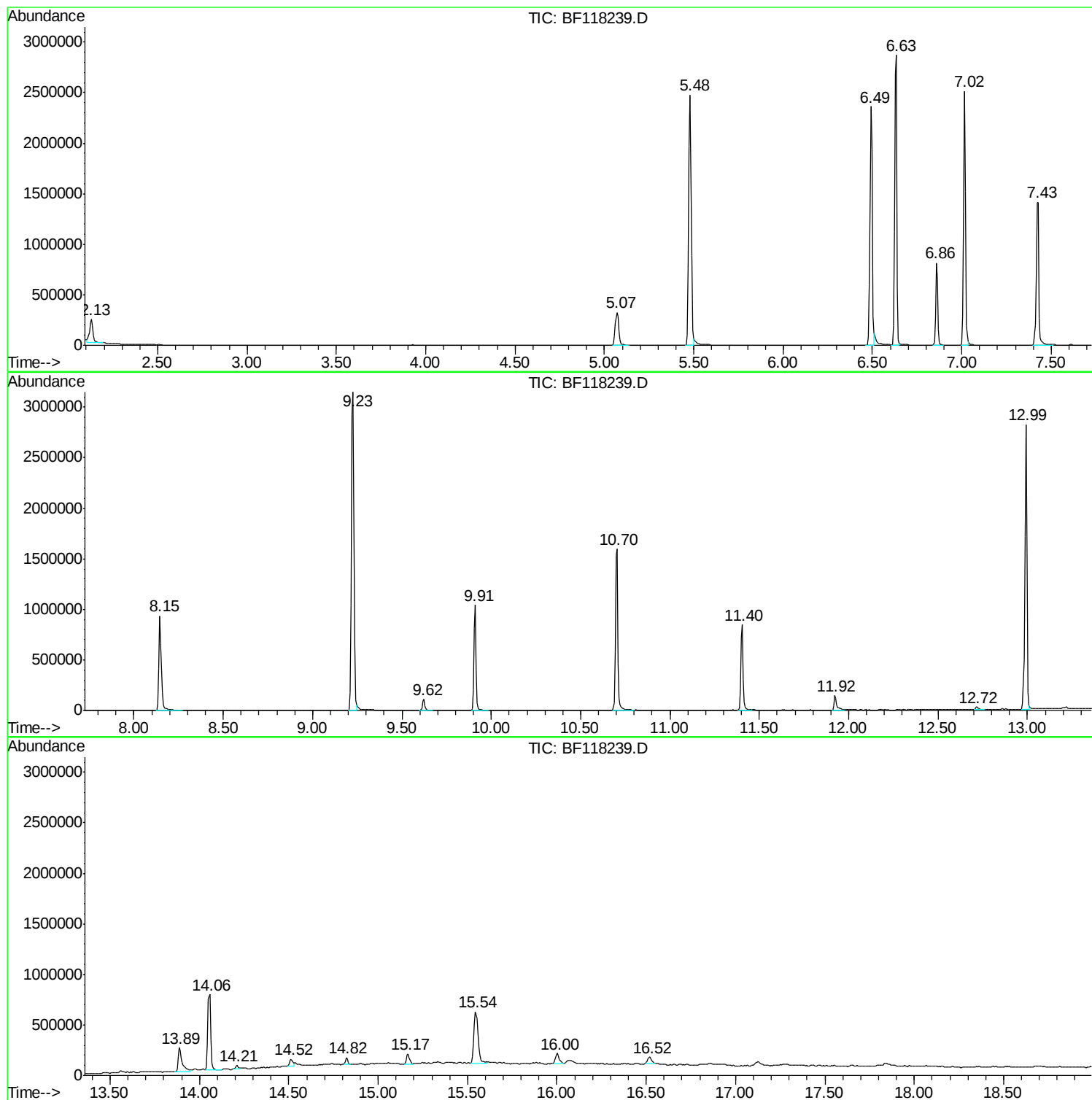
Sum of corrected areas: 24583209

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121719\
Data File : BF118239.D
Acq On : 17 Dec 2019 21:16
Operator : JU
Sample : K6324-05
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
EP-13

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF120619.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121719\
Data File : BF118239.D
Acq On : 17 Dec 2019 21:16
Operator : JU
Sample : K6324-05
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
EP-13

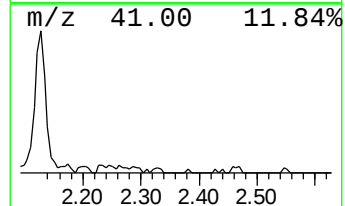
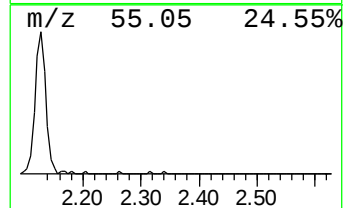
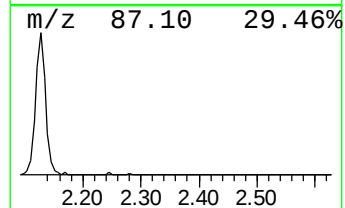
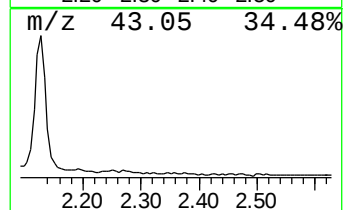
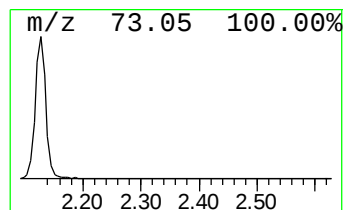
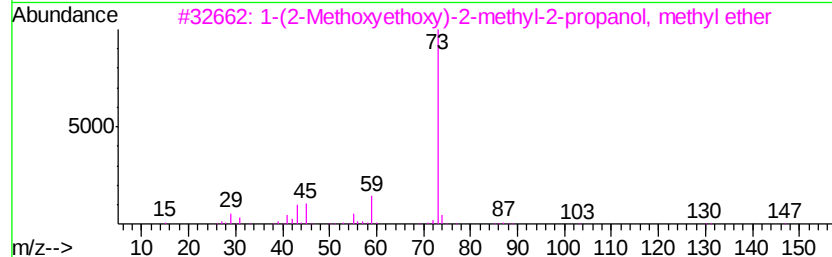
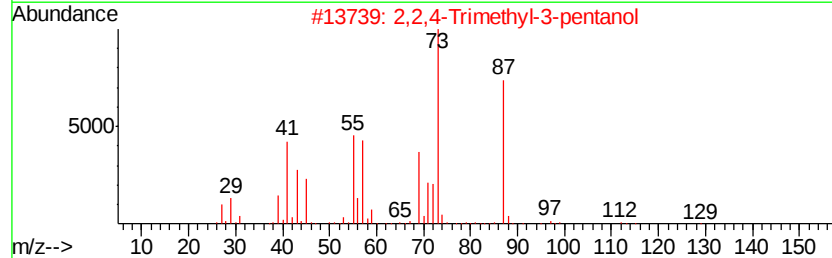
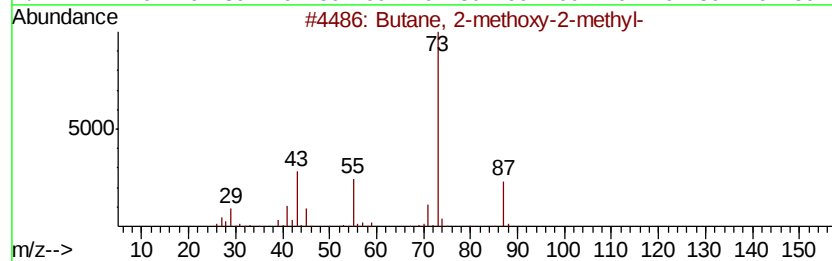
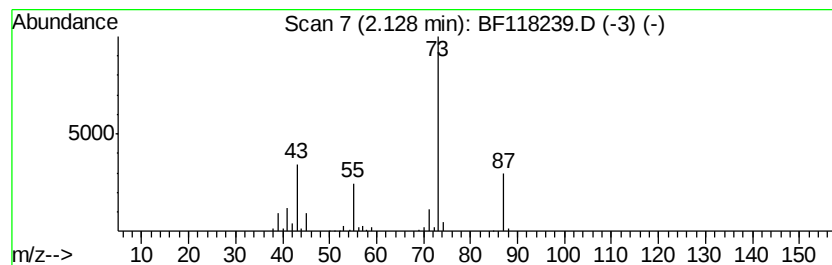
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF120619.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.13	8.98 ng	310234	1,4-Dichlorobenzene-d4	6.86

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2		2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	40
3		1-(2-Methoxyethoxy)-2-methyl-2-propanol	162	C8H18O3	1000364-24-4	28
4		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	17
5		Silane, tetramethyl-	88	C4H12Si	000075-76-3	12



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121719\
Data File : BF118239.D
Acq On : 17 Dec 2019 21:16
Operator : JU
Sample : K6324-05
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
EP-13

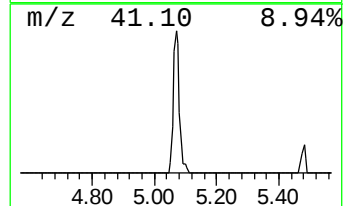
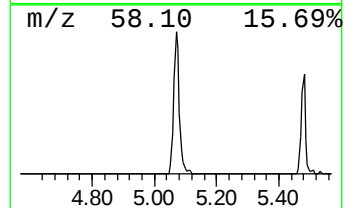
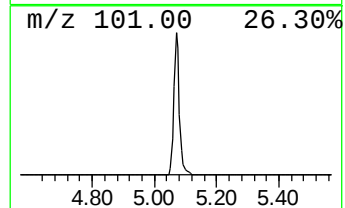
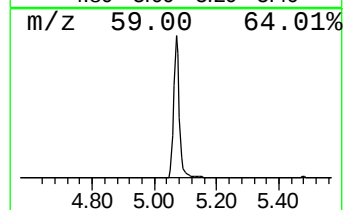
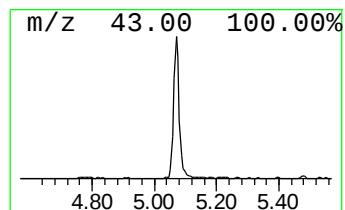
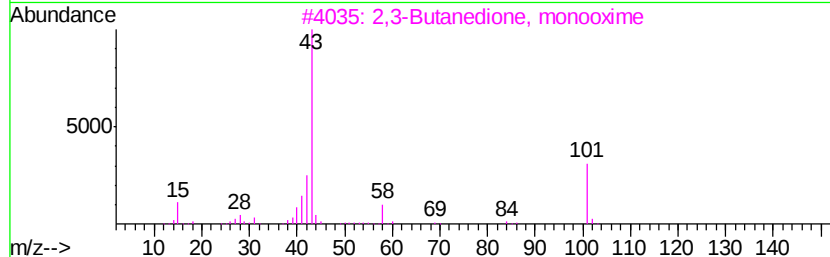
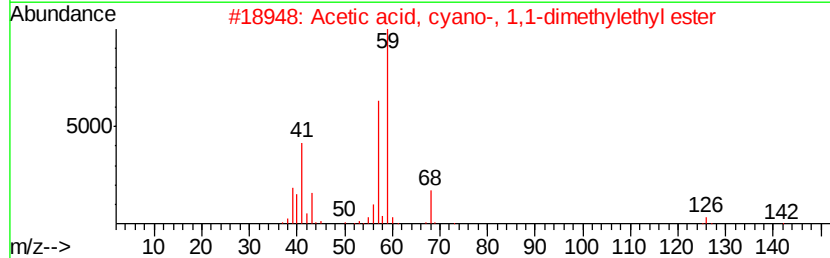
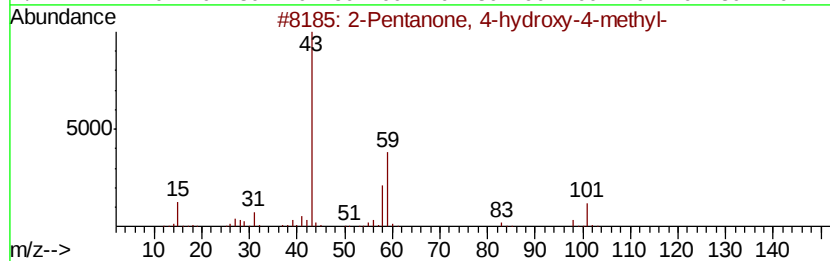
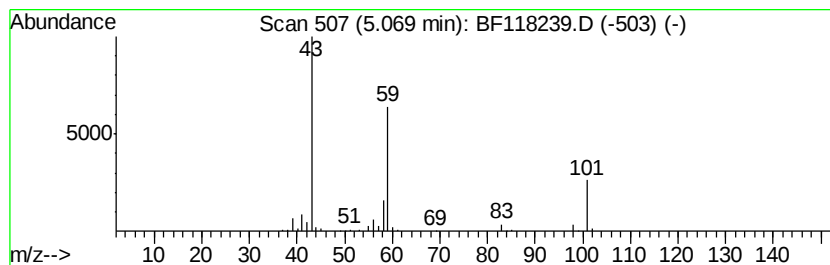
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF120619.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.07	11.92 ng	411989	1,4-Dichlorobenzene-d4	6.86

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2		Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	25
3		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	25
4		Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9
5		5-Hexen-2-one	98	C6H10O	000109-49-9	9



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121719\
Data File : BF118239.D
Acq On : 17 Dec 2019 21:16
Operator : JU
Sample : K6324-05
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
EP-13

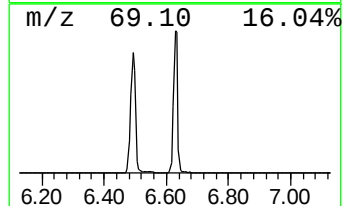
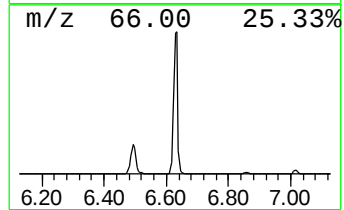
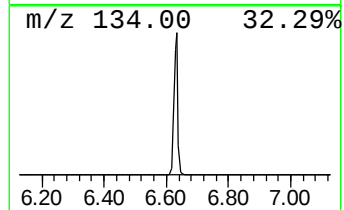
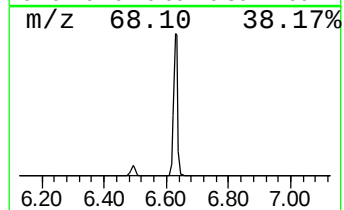
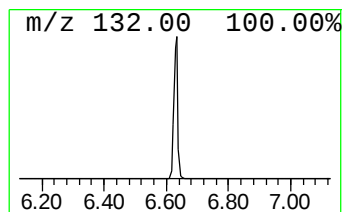
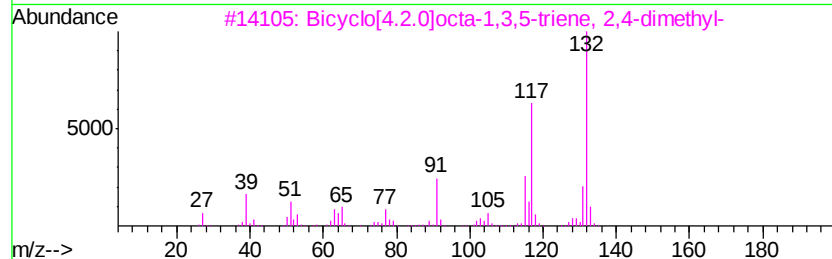
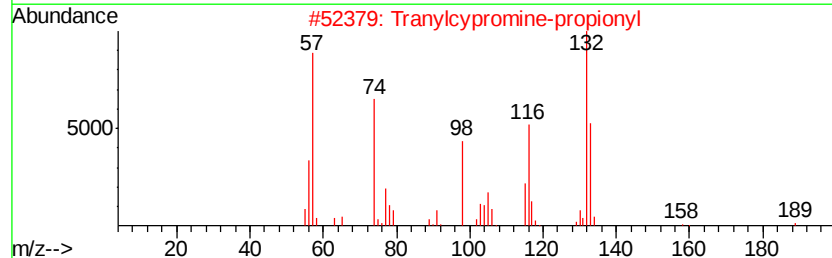
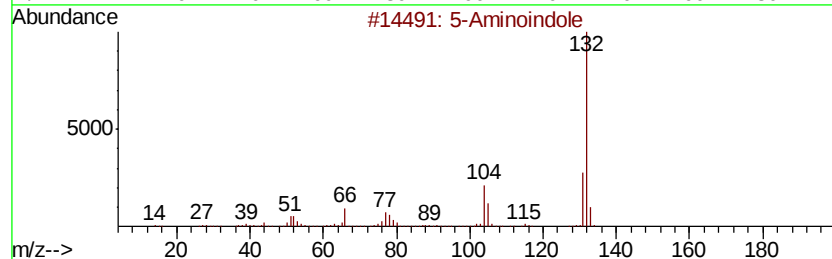
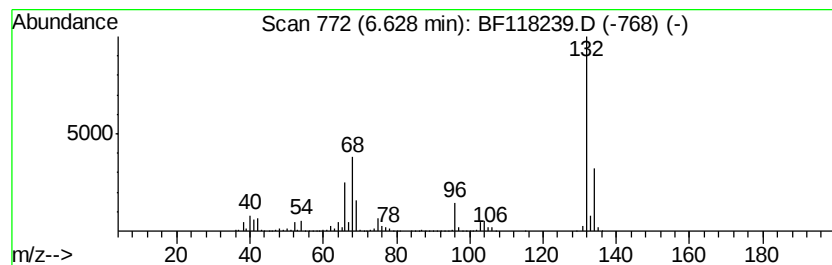
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF120619.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 3 unknown6.63 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.63	76.58 ng	2646840	1,4-Dichlorobenzene-d4	6.86

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Aminoindole	132	C8H8N2	005192-03-0	27
2		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	16
3		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
4		2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9
5		Xenon	132	Xe	007440-63-3	9



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121719\
Data File : BF118239.D
Acq On : 17 Dec 2019 21:16
Operator : JU
Sample : K6324-05
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
EP-13

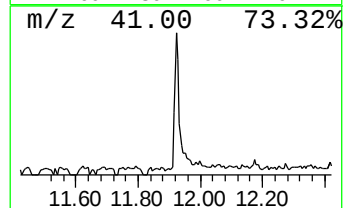
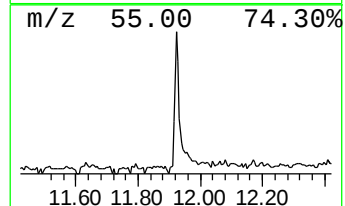
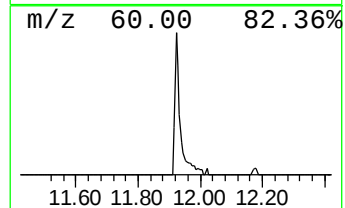
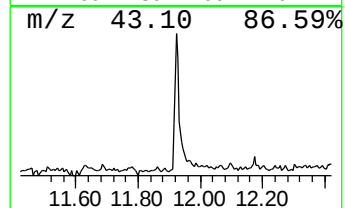
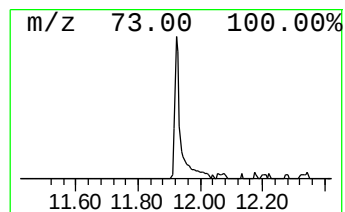
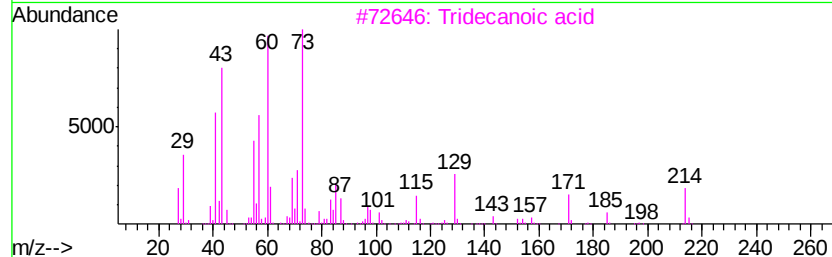
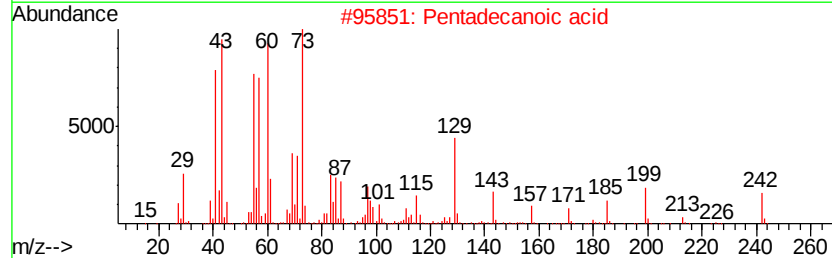
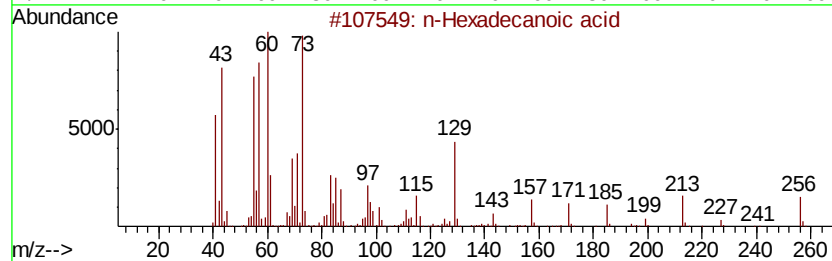
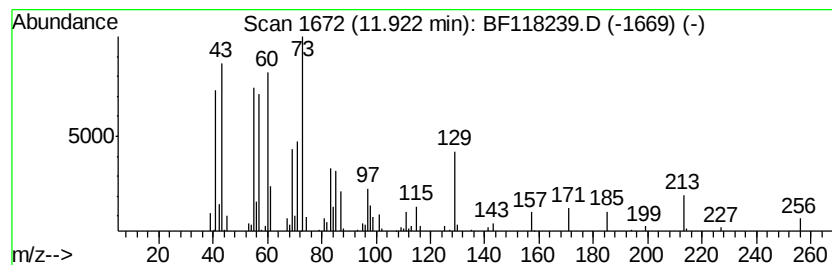
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF120619.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 n-Hexadecanoic acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.92	4.22 ng	168333	Phenanthrene-d10	11.40

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	96
2		Pentadecanoic acid	242	C15H30O2	001002-84-2	83
3		Tridecanoic acid	214	C13H26O2	000638-53-9	68
4		Sulfurous acid, octadecyl 2-prop...	376	C21H44O3S	1000309-12-7	18
5		Oxalic acid, isobutyl tetradecyl...	342	C20H38O4	1000309-37-9	11



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121719\
Data File : BF118239.D
Acq On : 17 Dec 2019 21:16
Operator : JU
Sample : K6324-05
Misc :
ALS Vial : 16 Sample Multiplier: 1

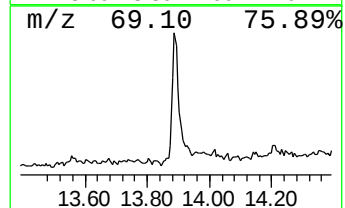
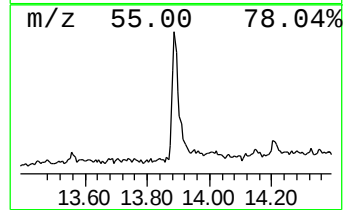
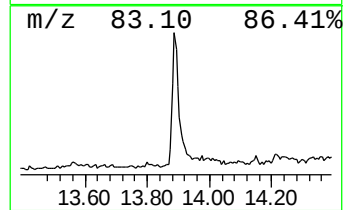
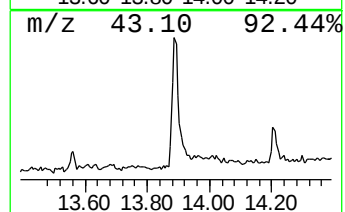
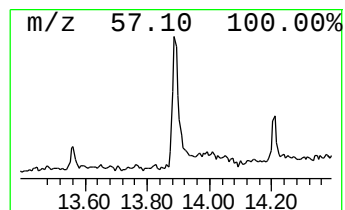
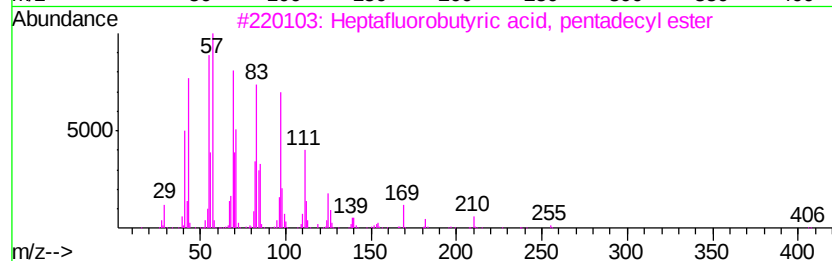
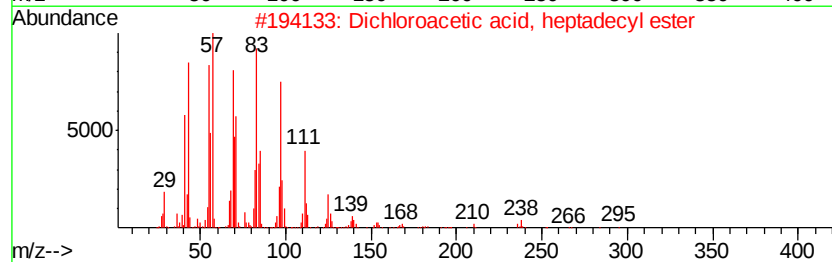
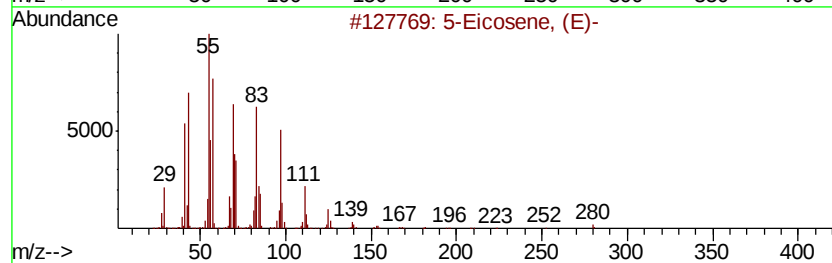
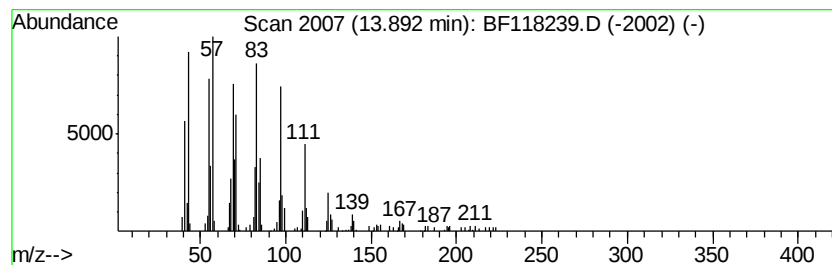
Instrument :
BNA_F
ClientSampleId :
EP-13

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF120619.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 5-Eicosene, (E)- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.		
13.89	9.93 ng	381694	Chrysene-d12	14.06		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	5-Eicosene, (E)-		280	C20H40	074685-30-6	98
2	Dichloroacetic acid, heptadecyl ...		366	C19H36Cl2O2	1000282-98-2	95
3	Heptafluorobutyric acid, pentade...		424	C19H31F7O2	959261-23-5	94
4	Pentafluoropropionic acid, penta...		374	C18H31F5O2	959092-08-1	94
5	Heptadecyl heptafluorobutyrate		452	C21H35F7O2	959085-66-6	94



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121719\
Data File : BF118239.D
Acq On : 17 Dec 2019 21:16
Operator : JU
Sample : K6324-05
Misc :
ALS Vial : 16 Sample Multiplier: 1

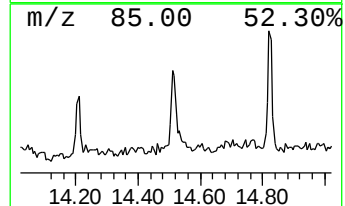
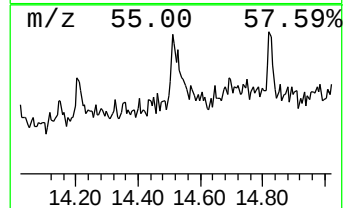
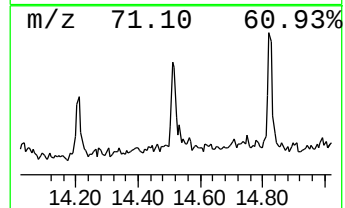
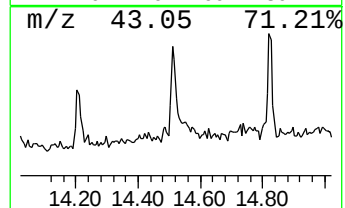
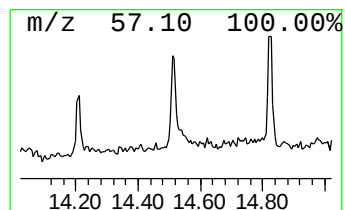
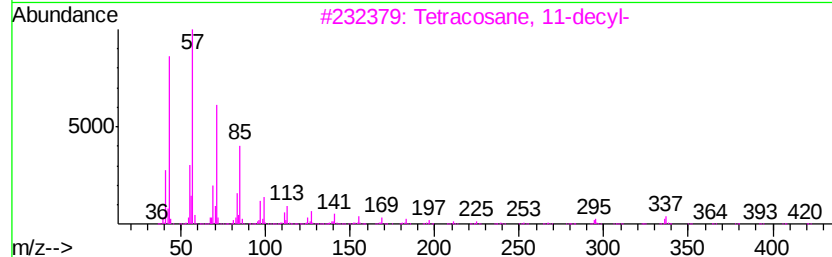
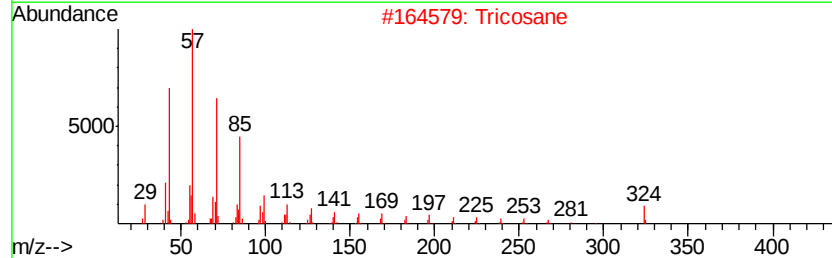
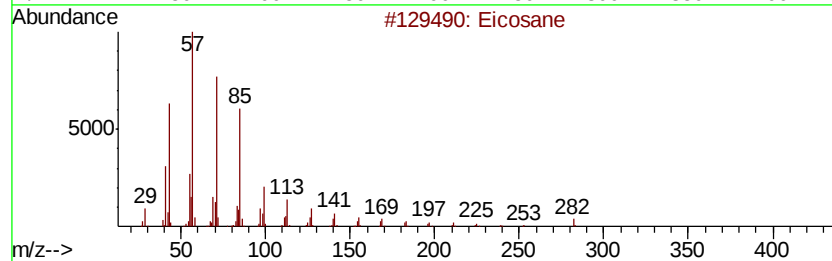
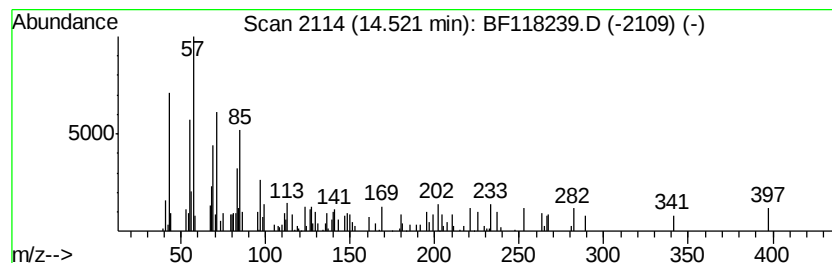
Instrument :
BNA_F
ClientSampleId :
EP-13

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF120619.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Eicosane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.52	2.13 ng	81724	Chrysene-d12	14.06
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Eicosane	282 C20H42	000112-95-8	93
2	Tricosane	324 C23H48	000638-67-5	90
3	Tetracosane, 11-decyl-	479 C34H70	055429-84-0	76
4	Hentriacontane	437 C31H64	000630-04-6	76
5	Nonane, 5-butyl-	184 C13H28	017312-63-9	70



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121719\
Data File : BF118239.D
Acq On : 17 Dec 2019 21:16
Operator : JU
Sample : K6324-05
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
EP-13

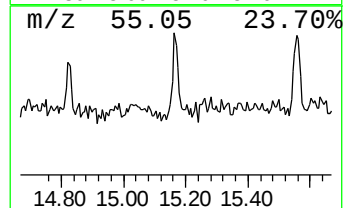
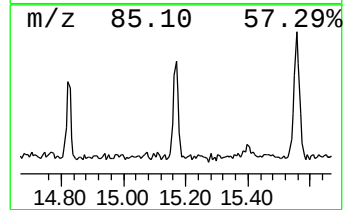
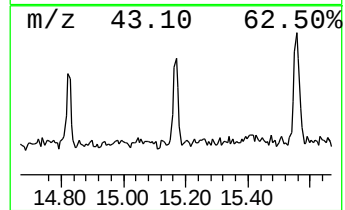
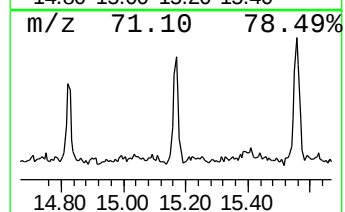
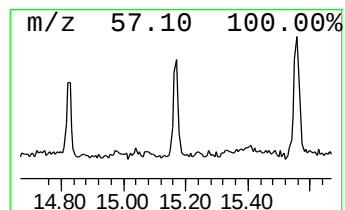
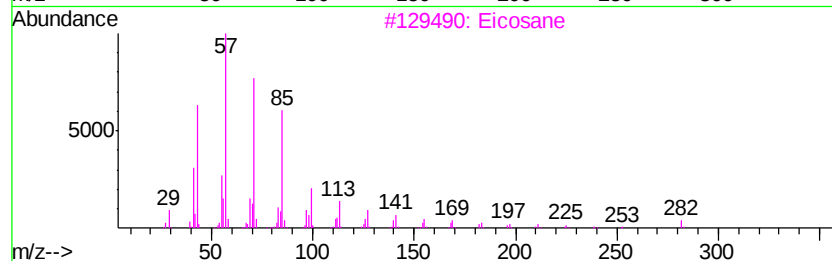
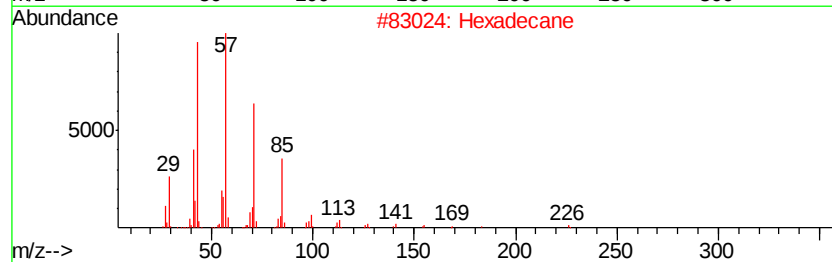
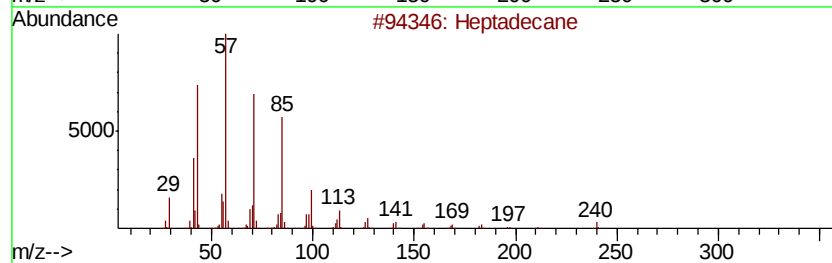
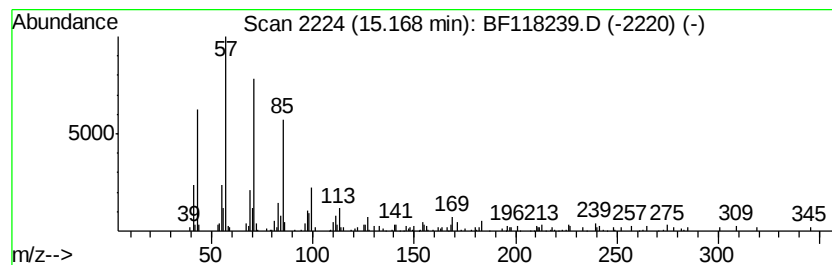
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF120619.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Heptadecane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.17	2.90 ng	115054	Perylene-d12	15.54

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane		240	C17H36	000629-78-7	94
2	Hexadecane		226	C16H34	000544-76-3	93
3	Eicosane		282	C20H42	000112-95-8	91
4	Heneicosane		296	C21H44	000629-94-7	91
5	Octadecane		254	C18H38	000593-45-3	86



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121719\
Data File : BF118239.D
Acq On : 17 Dec 2019 21:16
Operator : JU
Sample : K6324-05
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
EP-13

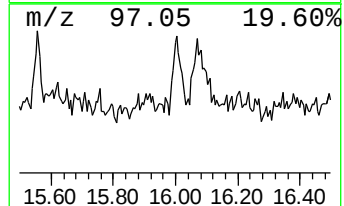
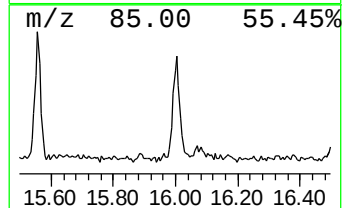
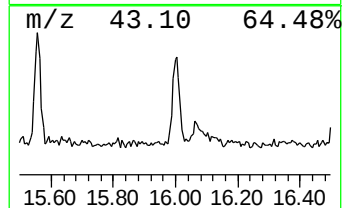
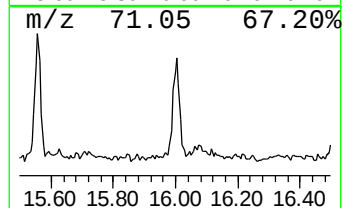
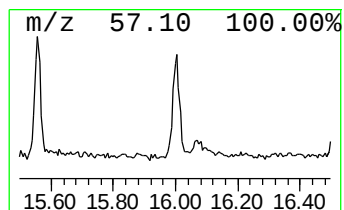
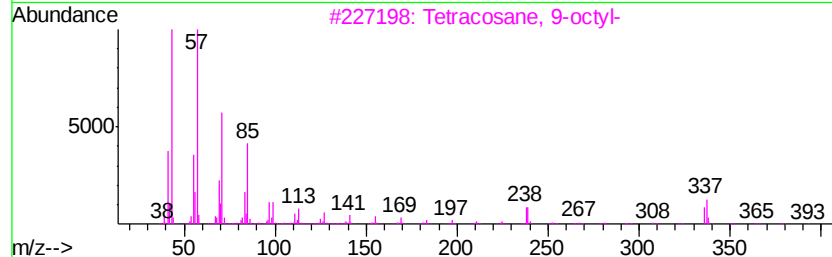
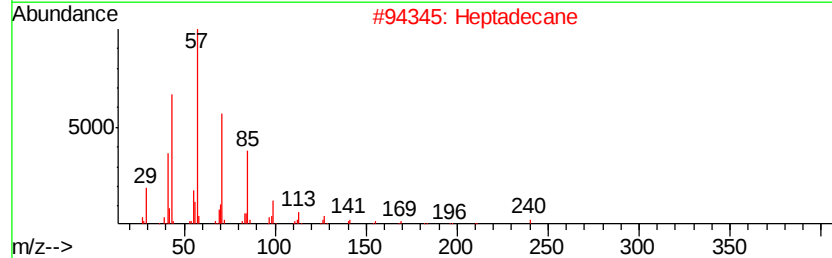
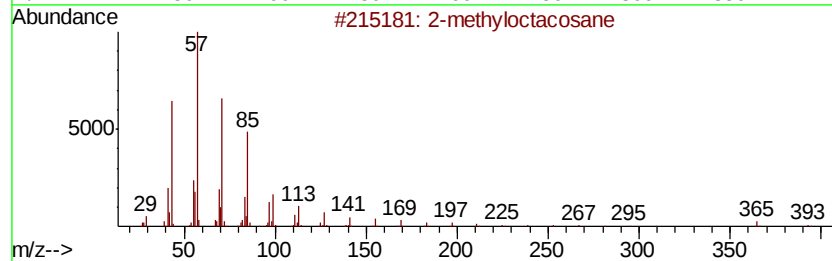
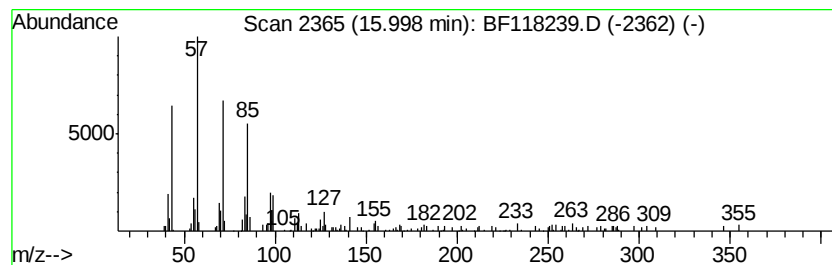
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF120619.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 2-methyloctacosane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.00	3.65 ng	144955	Perylene-d12	15.54

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-methyloctacosane		408	C29H60	1000376-72-8	91
2	Heptadecane		240	C17H36	000629-78-7	90
3	Tetracosane, 9-octyl-		451	C32H66	055401-54-2	86
4	Eicosane		282	C20H42	000112-95-8	81
5	Hexadecane, 2-methyl-		240	C17H36	001560-92-5	76



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121719\
Data File : BF118239.D
Acq On : 17 Dec 2019 21:16
Operator : JU
Sample : K6324-05
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
EP-13

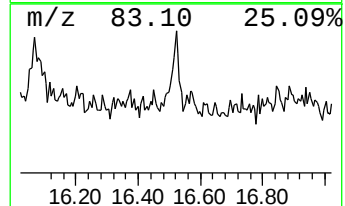
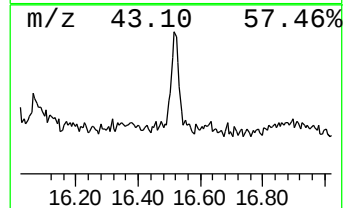
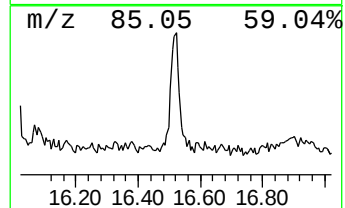
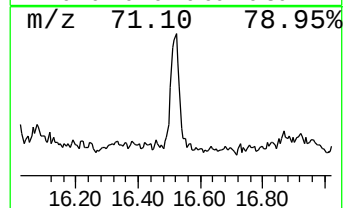
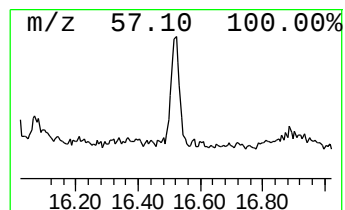
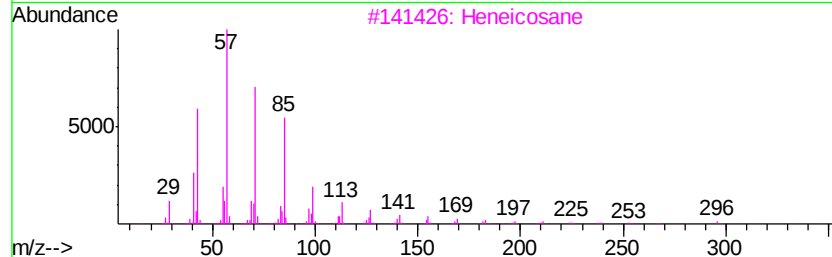
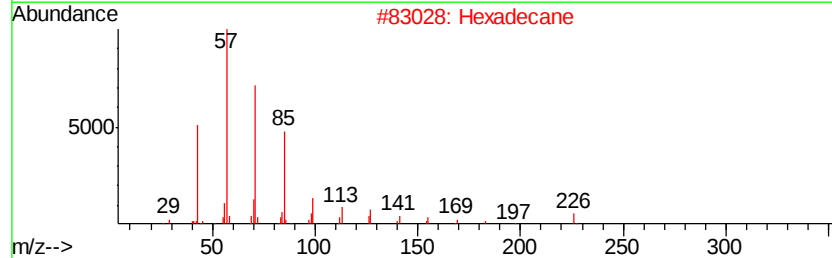
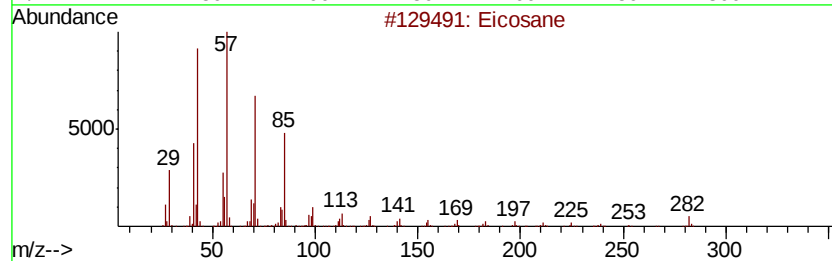
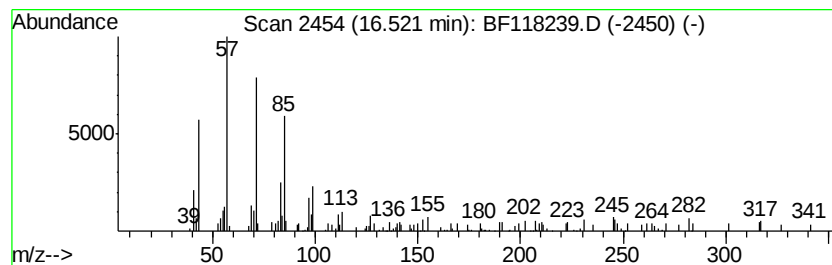
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF120619.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Hexadecane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.52	2.47 ng	98012	Perylene-d12	15.54

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane	282	C20H42	000112-95-8	89
2		Hexadecane	226	C16H34	000544-76-3	76
3		Heneicosane	296	C21H44	000629-94-7	74
4		2-methyloctacosane	408	C29H60	1000376-72-8	74
5		Heptadecane	240	C17H36	000629-78-7	72



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF121719\
Data File : BF118239.D
Acq On : 17 Dec 2019 21:16
Operator : JU
Sample : K6324-05
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
EP-13

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF120619.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
Butane, 2-methoxy...	2.13	9.0	ng	310234	1	6.86	691219	20.0
2-Pentanone, 4-hy...	5.07	11.9	ng	411989	1	6.86	691219	20.0
unknown6.63	6.63	76.6	ng	2646840	1	6.86	691219	20.0
n-Hexadecanoic acid	11.92	4.2	ng	168333	4	11.40	797330	20.0
5-Eicosene, (E)-	13.89	9.9	ng	381694	5	14.06	769120	20.0
Eicosane	14.52	2.1	ng	81724	5	14.06	769120	20.0
Heptadecane	15.17	2.9	ng	115054	6	15.54	793377	20.0
2-methyloctacosane	16.00	3.6	ng	144955	6	15.54	793377	20.0
Hexadecane	16.52	2.5	ng	98012	6	15.54	793377	20.0