

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF121719\
 Data File : BF118238.D
 Acq On : 17 Dec 2019 20:48
 Operator : JU
 Sample : K6329-01
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 121619-CA-COMP

Integration Parameters: rteint.p

Integrator: RTE

Smoothing : ON

Sampling : 1

Start Thrs: 0.2

Stop Thrs : 0

Filtering: 5

Min Area: 3 % of largest Peak

Max Peaks: 100

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-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	2.122	3	6	21	rVB	295715	373528	11.40%	1.238%
2	4.916	473	481	487	rVB	29832	38266	1.17%	0.127%
3	5.069	503	507	515	rBV	421987	524852	16.02%	1.740%
4	5.269	535	541	544	rBV2	40011	46044	1.41%	0.153%
5	5.481	573	577	580	rBV	2716876	2457649	75.01%	8.146%
6	5.628	599	602	607	rBV3	35149	50101	1.53%	0.166%
7	5.716	612	617	625	rVB2	28779	41103	1.25%	0.136%
8	5.881	643	645	648	rVB	43015	38580	1.18%	0.128%
9	6.016	663	668	672	rVB3	50236	64679	1.97%	0.214%
10	6.104	678	683	690	rBV2	108071	147823	4.51%	0.490%
11	6.298	710	716	718	rBV4	39681	57375	1.75%	0.190%
12	6.351	723	725	728	rVB	44118	39077	1.19%	0.130%
13	6.393	728	732	737	rBV3	58311	87709	2.68%	0.291%
14	6.498	744	750	758	rBV	2553638	2796318	85.34%	9.269%
15	6.634	768	773	776	rBV	3186923	2901814	88.56%	9.618%
16	6.857	808	811	818	rVB	875681	778400	23.76%	2.580%
17	7.016	833	838	841	rBV	2817601	2335035	71.26%	7.740%
18	7.134	855	858	861	rVB2	44732	45033	1.37%	0.149%
19	7.257	873	879	881	rBV4	28827	38951	1.19%	0.129%
20	7.422	903	907	916	rVB	1631695	1693336	51.68%	5.613%
21	7.504	916	921	924	rVB5	33937	50559	1.54%	0.168%
22	8.145	1025	1030	1036	rBV	1080159	934622	28.52%	3.098%
23	8.204	1036	1040	1043	rVB2	28155	41028	1.25%	0.136%
24	9.228	1209	1214	1217	rBV	3940841	3276563	100.00%	10.860%
25	9.622	1277	1281	1290	rBV	148154	143889	4.39%	0.477%
26	9.910	1326	1330	1338	rBV	1208351	1039888	31.74%	3.447%
27	10.704	1461	1465	1475	rBV	2174232	1987988	60.67%	6.589%
28	11.404	1580	1584	1587	rBV	1003853	911155	27.81%	3.020%
29	11.922	1667	1672	1680	rBV2	223740	268120	8.18%	0.889%
30	12.621	1788	1791	1796	rBV	97927	98348	3.00%	0.326%
31	12.716	1803	1807	1812	rBV2	56229	76358	2.33%	0.253%
32	12.851	1825	1830	1834	rBV	88559	106260	3.24%	0.352%
33	12.992	1850	1854	1857	rBV	3183694	2763463	84.34%	9.160%
34	13.216	1888	1892	1895	rVB3	41936	46678	1.42%	0.155%

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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

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

35	13.563	1947	1951	1954	rBV	55197	66412	2.03%	0.220%
36	13.886	2002	2006	2017	rBV	437149	574303	17.53%	1.904%
37	14.057	2030	2035	2038	rBV	872410	850483	25.96%	2.819%
38	14.210	2058	2061	2068	rBV	97047	105284	3.21%	0.349%
39	14.515	2110	2113	2121	rVB2	140686	216112	6.60%	0.716%
40	14.821	2162	2165	2169	rVB	139414	156963	4.79%	0.520%
41	14.904	2176	2179	2185	rVB2	54680	65808	2.01%	0.218%
42	15.168	2220	2224	2228	rVB	188455	202270	6.17%	0.670%
43	15.551	2284	2289	2299	rBV2	606524	961109	29.33%	3.186%
44	16.004	2361	2366	2372	rVB	179690	262838	8.02%	0.871%
45	16.068	2374	2377	2388	rVB7	47674	92894	2.84%	0.308%
46	16.521	2449	2454	2459	rBV2	102185	170900	5.22%	0.566%
47	17.127	2553	2557	2566	rVB4	66171	143820	4.39%	0.477%

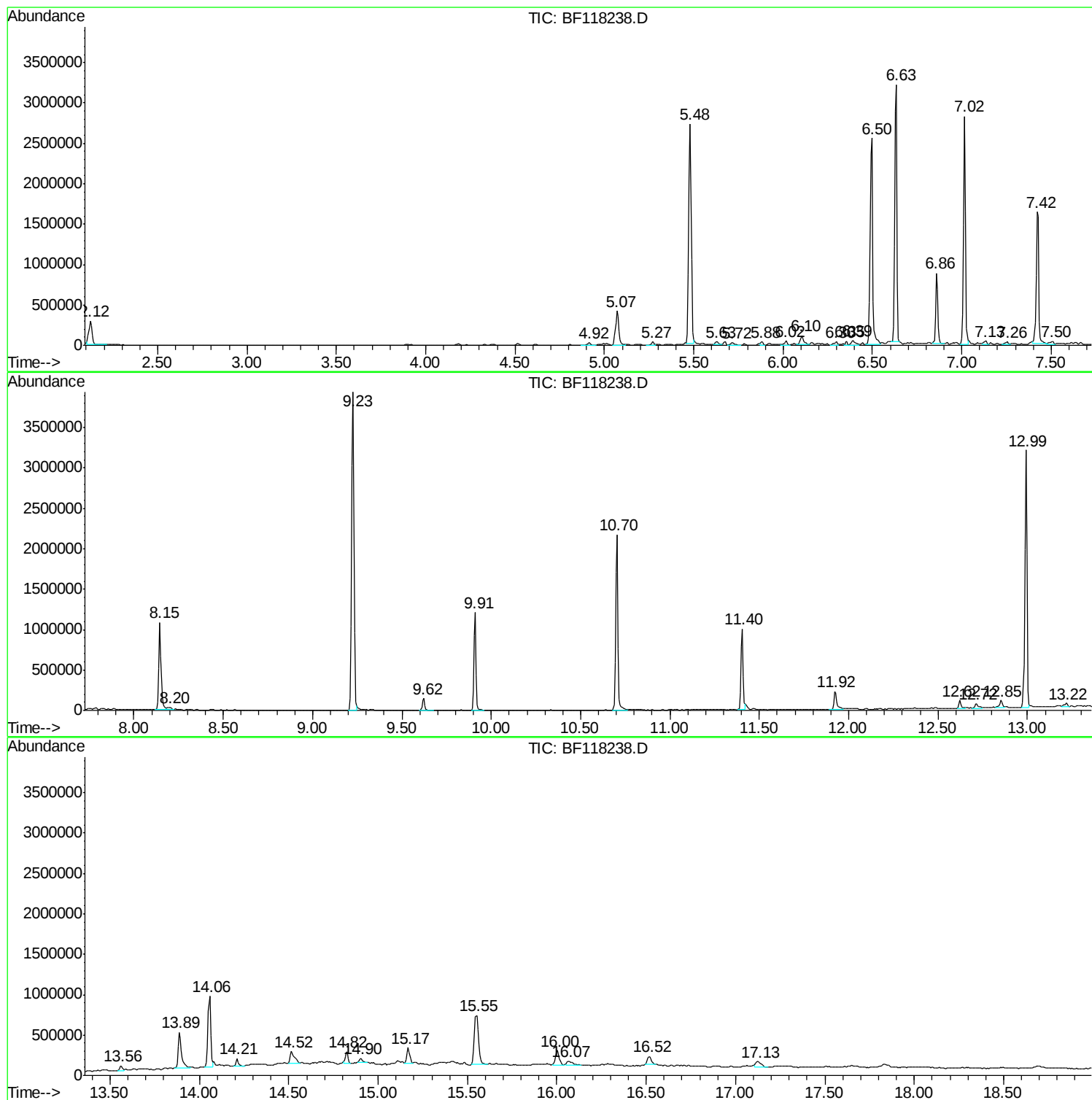
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF120619.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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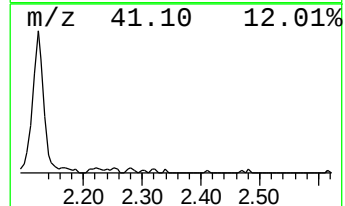
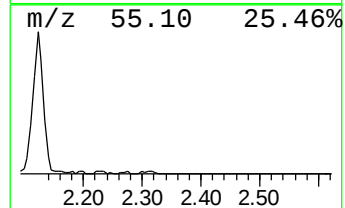
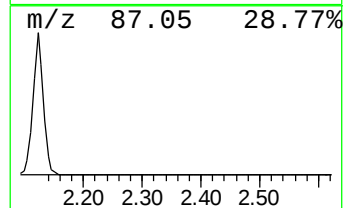
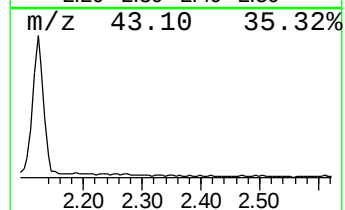
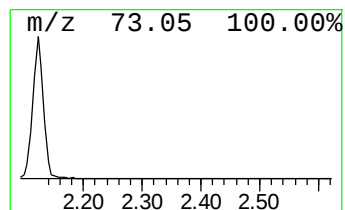
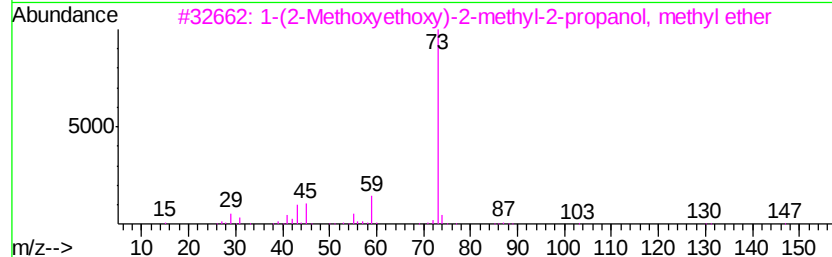
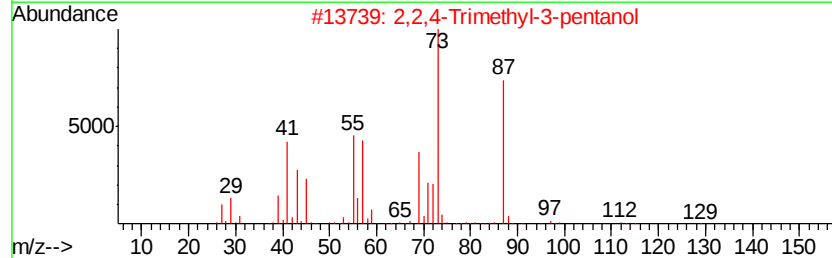
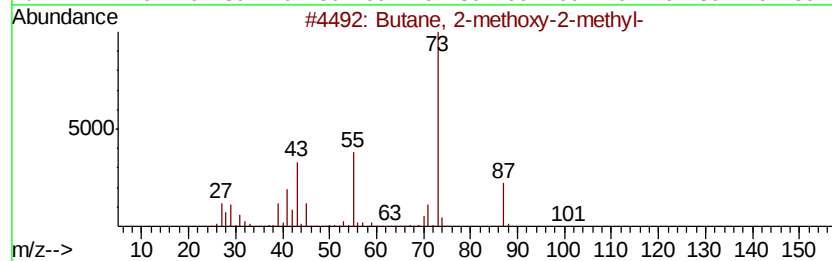
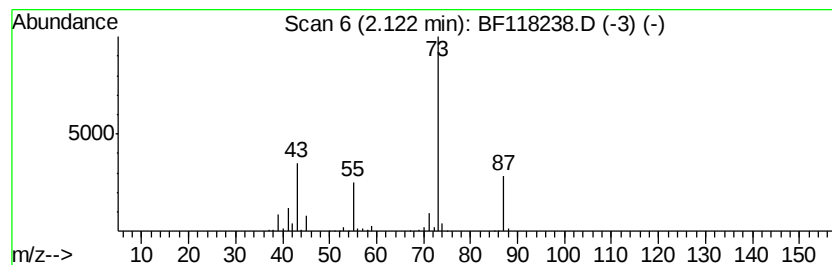
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF120619.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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.
2.12	9.60 ng	373528	1,4-Dichlorobenzene-d4	6.86

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	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-(2-Methoxyethoxy)-2-methyl-2-propanol	162	C8H18O3	1000364-24-4	25
4		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	23
5		Silane, tetramethyl-	88	C4H12Si	000075-76-3	17



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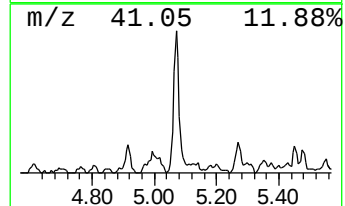
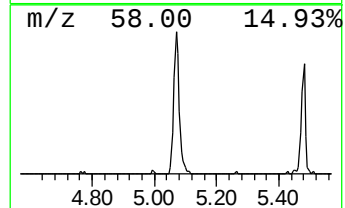
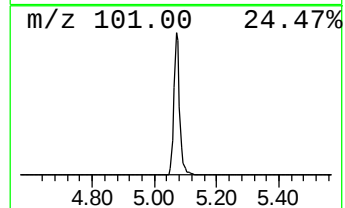
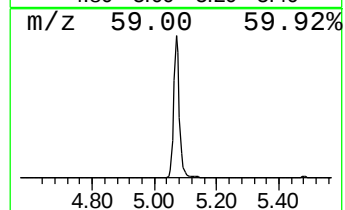
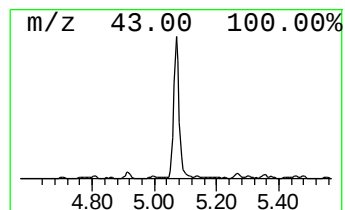
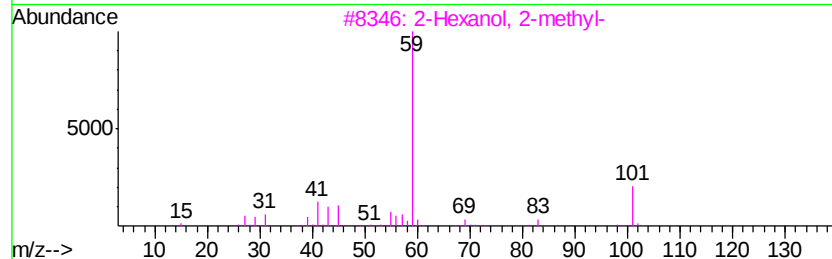
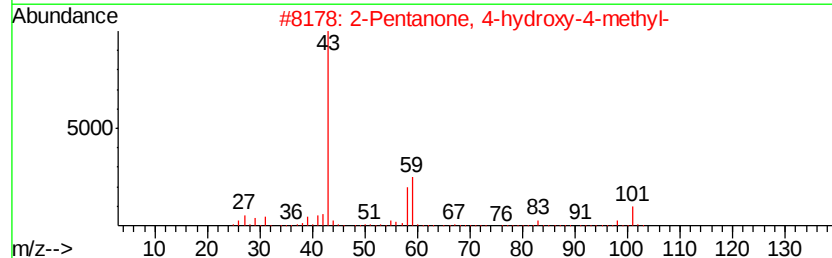
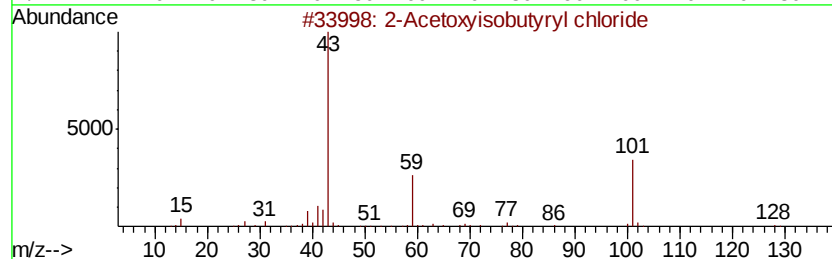
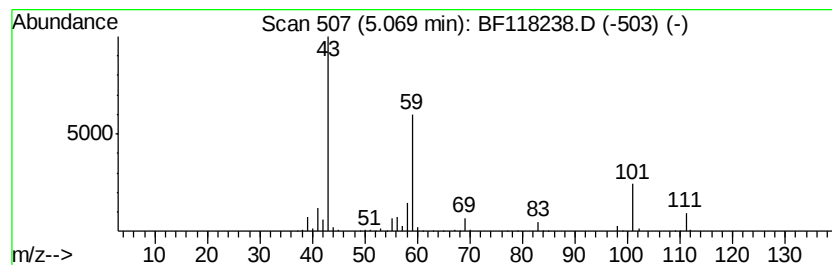
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF120619.M
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 2-Acetoxyisobutyryl chloride Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.07	13.49 ng	524852	1,4-Dichlorobenzene-d4	6.86

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Acetoxyisobutyryl chloride	164	C6H9ClO3	040635-66-3	59
2		2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	53
3		2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	38
4		Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	36
5		n-Propyl t-butyl ether	116	C7H16O	1000334-81-9	25



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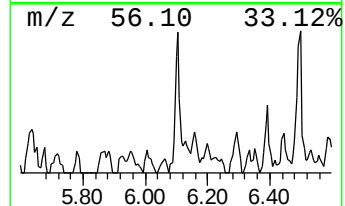
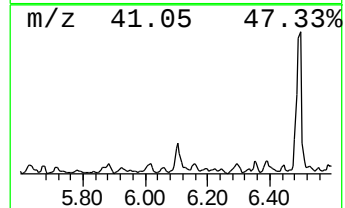
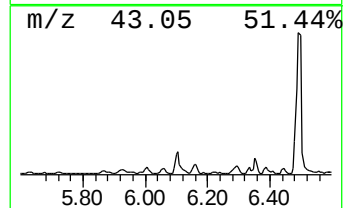
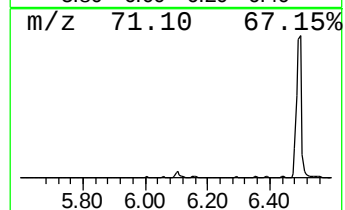
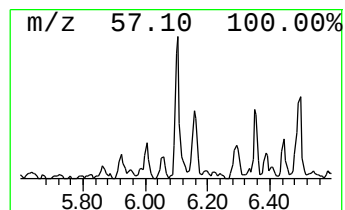
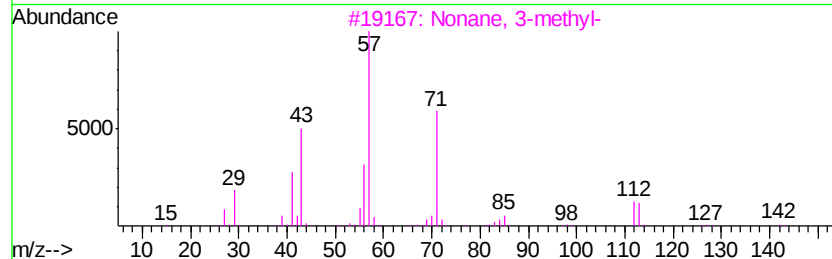
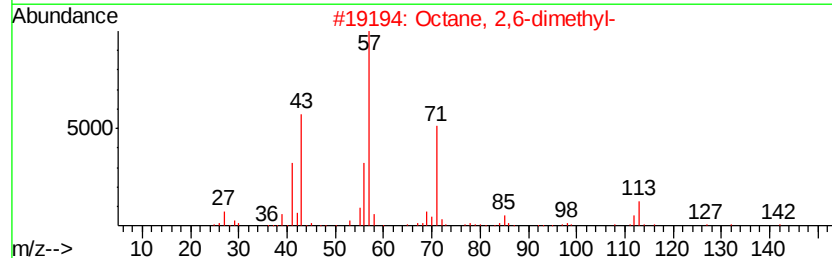
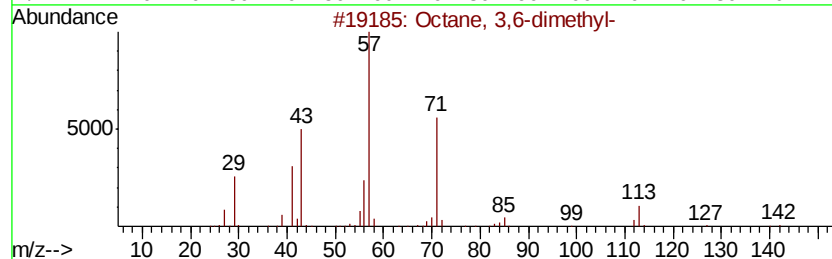
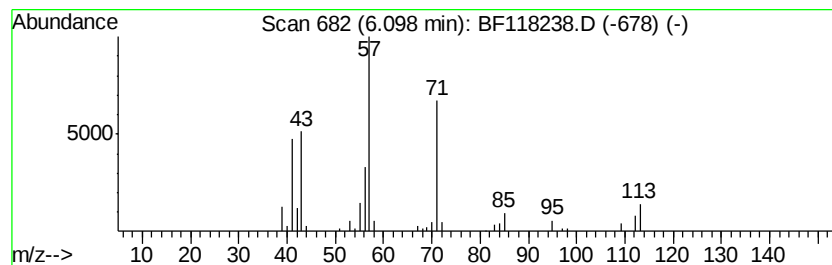
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 3 Octane, 3,6-dimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.10	3.80 ng	147823	1,4-Dichlorobenzene-d4	6.86

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octane, 3,6-dimethyl-	142	C10H22	015869-94-0	59
2		Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	59
3		Nonane, 3-methyl-	142	C10H22	005911-04-6	53
4		Undecane, 4,5-dimethyl-	184	C13H28	017312-79-7	50
5		Sulfurous acid, butyl 2-ethylhex...	250	C12H26O3S	1000309-18-8	50



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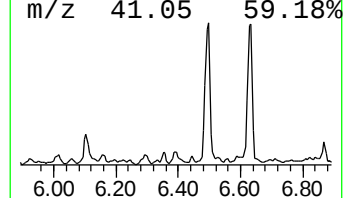
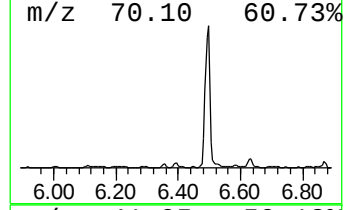
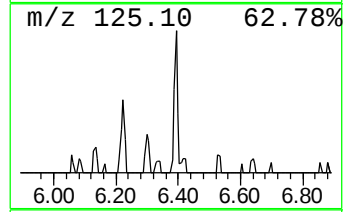
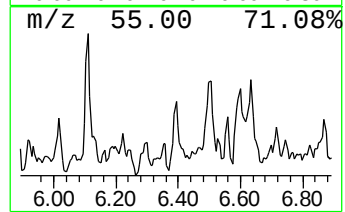
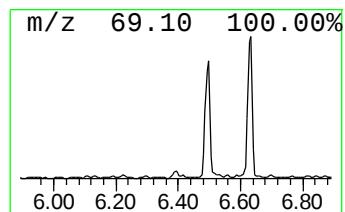
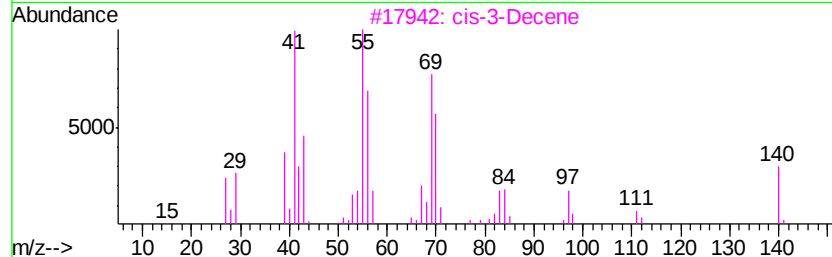
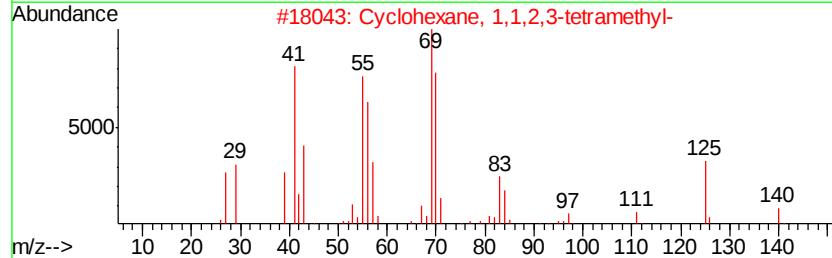
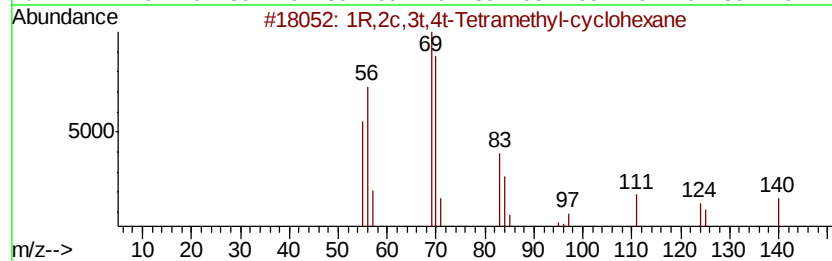
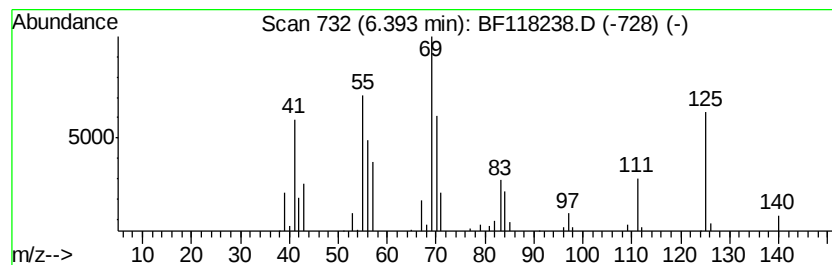
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 1R,2c,3t,4t-Tetramethyl-cyc... Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.39	2.25 ng	87709	1,4-Dichlorobenzene-d4	6.86

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1R,2c,3t,4t-Tetramethyl-cyclohexane	140	C10H20	1000144-07-3	81
2		Cyclohexane, 1,1,2,3-tetramethyl-	140	C10H20	006783-92-2	72
3		cis-3-Decene	140	C10H20	019398-86-8	62
4		1-Octene, 3,7-dimethyl-	140	C10H20	004984-01-4	49
5		Cyclopentane, butyl-	126	C9H18	002040-95-1	49



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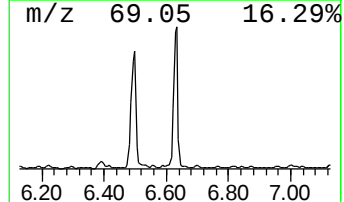
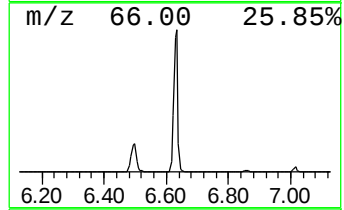
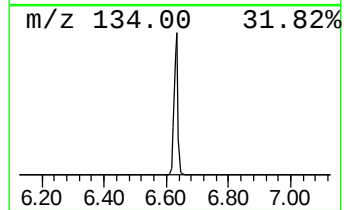
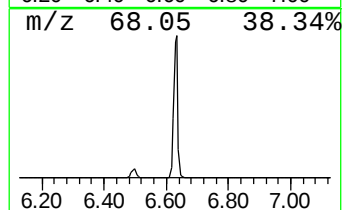
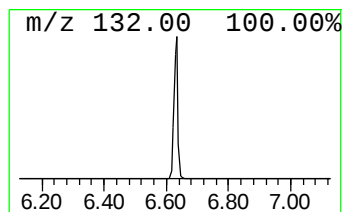
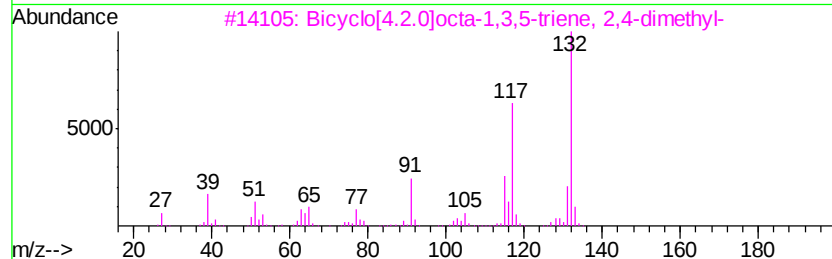
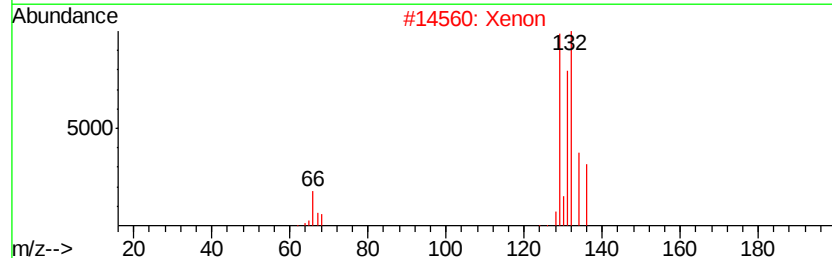
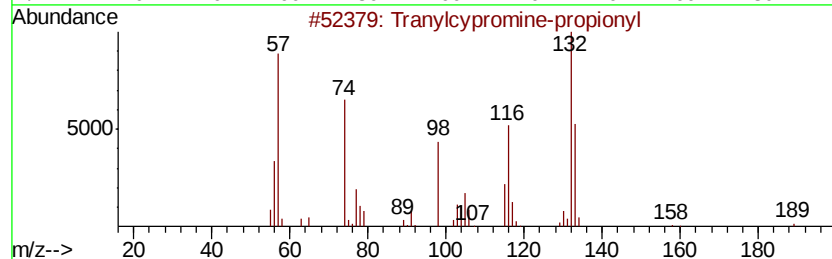
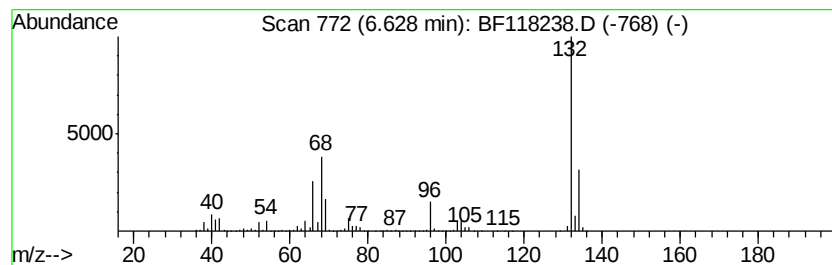
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 unknown6.63 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.63	74.56 ng	2901810	1,4-Dichlorobenzene-d4	6.86

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tranvlcvpromine-propionyl	189	C12H15NO	1000123-86-3	16
2		Xenon	132	Xe	007440-63-3	9
3		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
4		Cyclopropanamine, 1-phenyl-	133	C9H11N	1000351-26-2	9
5		1-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-59-9	9



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121719\
Data File : BF118238.D
Acq On : 17 Dec 2019 20:48
Operator : JU
Sample : K6329-01
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
121619-CA-COMP

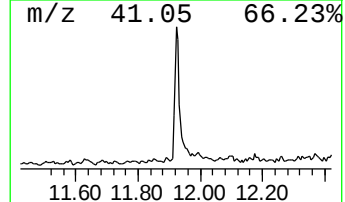
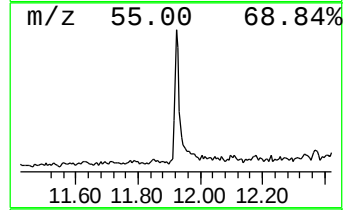
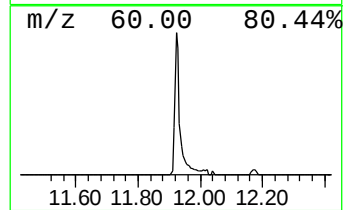
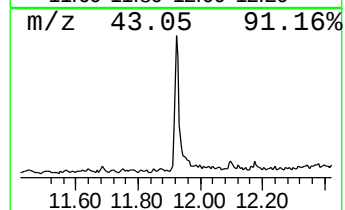
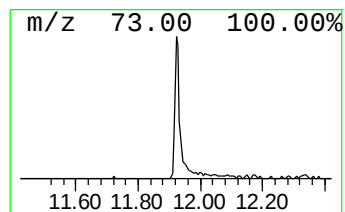
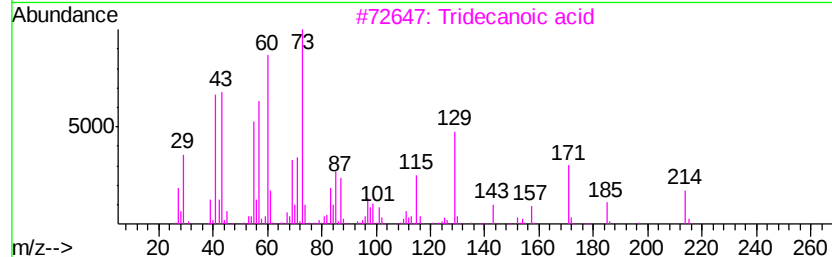
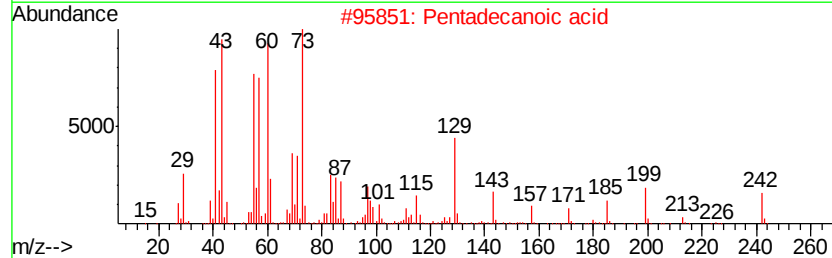
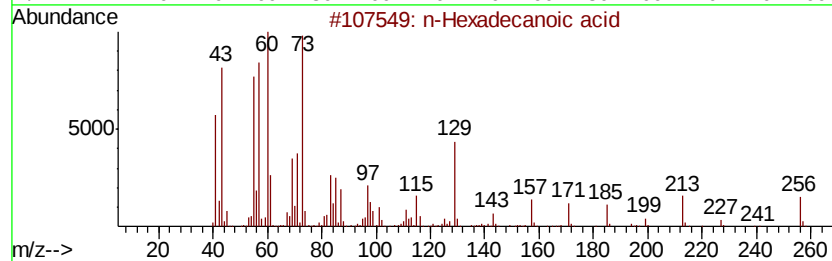
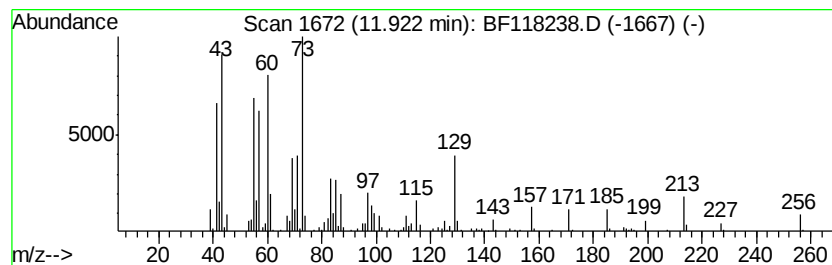
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 n-Hexadecanoic acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.92	5.89 ng	268120	Phenanthrene-d10	11.40

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2	Pentadecanoic acid	242	C15H30O2	001002-84-2	94
3	Tridecanoic acid	214	C13H26O2	000638-53-9	93
4	Tetradecanoic acid	228	C14H28O2	000544-63-8	93
5	n-Decanoic acid	172	C10H20O2	000334-48-5	81



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121719\
Data File : BF118238.D
Acq On : 17 Dec 2019 20:48
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Misc :
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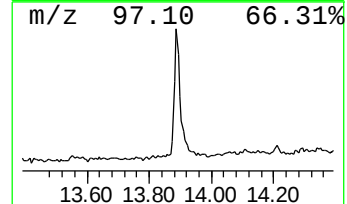
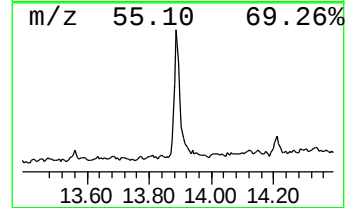
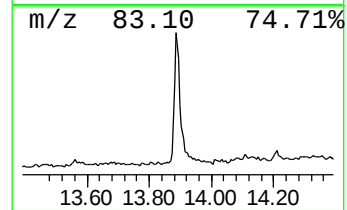
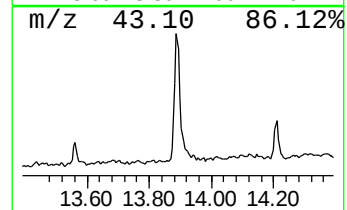
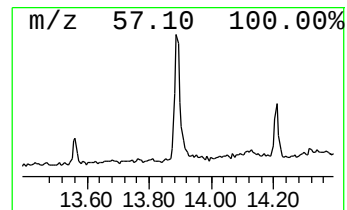
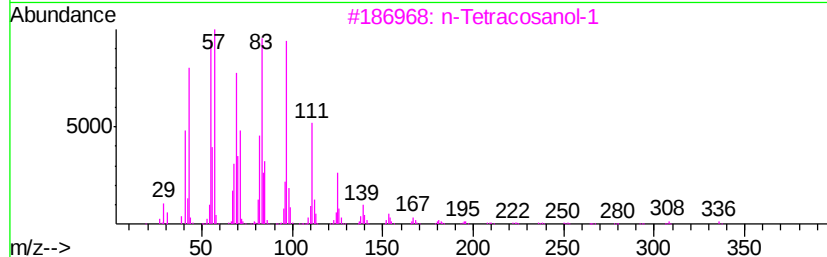
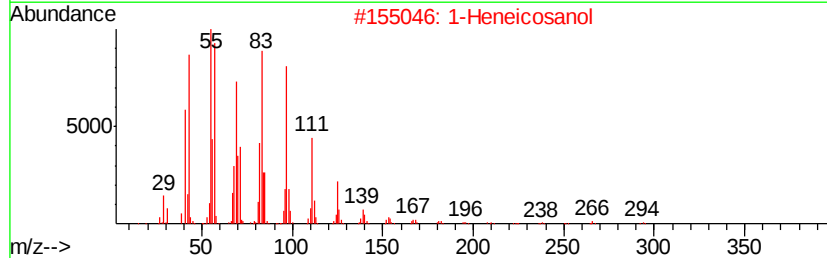
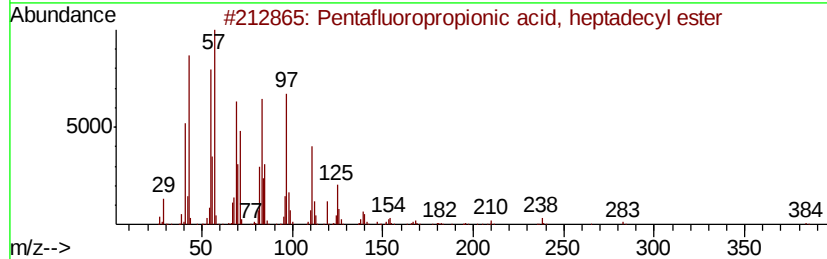
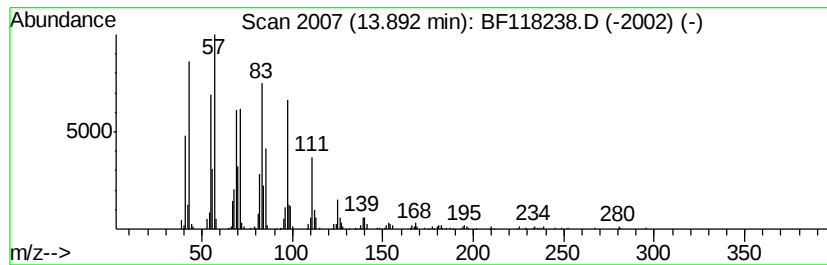
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Pentafluoropropionic acid, ... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.89	13.51 ng	574303	Chrysene-d12	14.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentafluoropropionic acid, hepta...	402	C20H35F5O2	959218-78-1	94
2		1-Heneicosanol	312	C21H44O	015594-90-8	93
3		n-Tetracosanol-1	354	C24H50O	000506-51-4	93
4		Behenic alcohol	326	C22H46O	000661-19-8	93
5		Nonadecyl pentafluoropropionate	430	C22H39F5O2	1000351-88-8	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121719\
Data File : BF118238.D
Acq On : 17 Dec 2019 20:48
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Misc :
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ClientSampleId :
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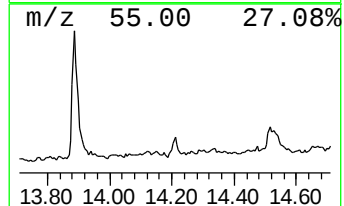
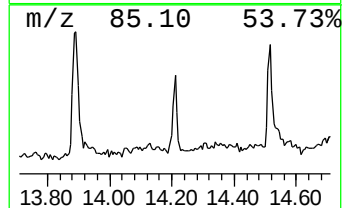
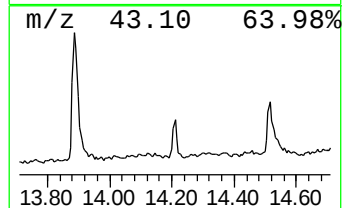
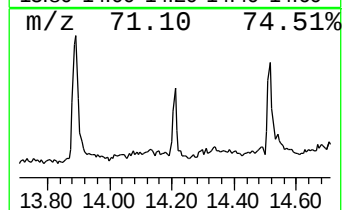
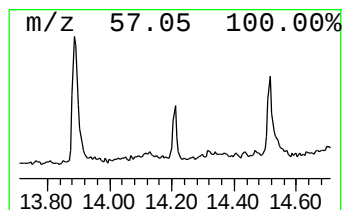
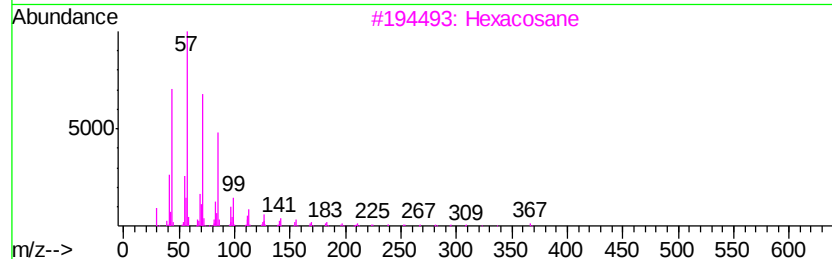
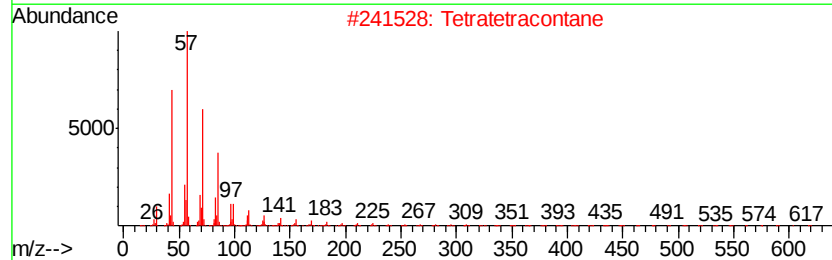
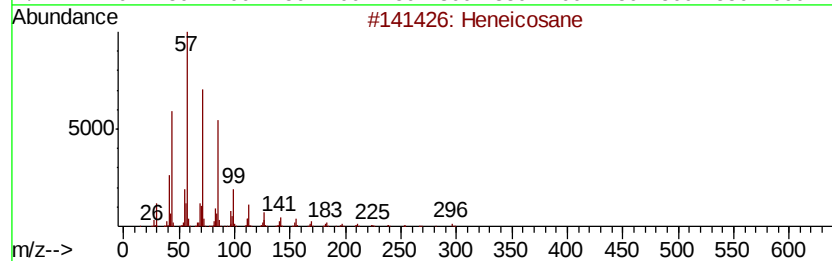
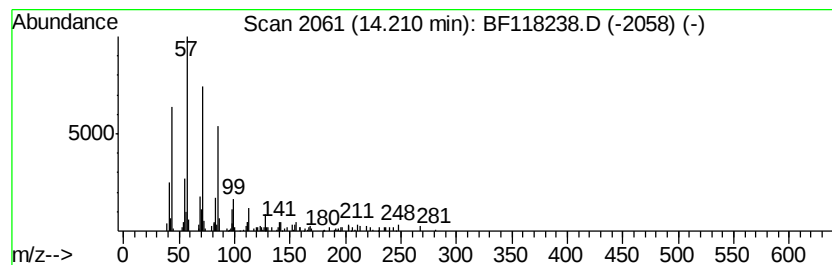
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Heneicosane Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.21	2.48 ng	105284	Chrysene-d12	14.06

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heneicosane	296	C21H44	000629-94-7	91
2			Tetratetracontane	619	C44H90	007098-22-8	91
3			Hexacosane	366	C26H54	000630-01-3	91
4			2-methyloctacosane	408	C29H60	1000376-72-8	90
5			Dodecane, 1-iodo-	296	C12H25I	004292-19-7	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121719\
Data File : BF118238.D
Acq On : 17 Dec 2019 20:48
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Instrument :
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ClientSampleId :
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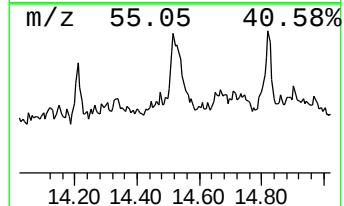
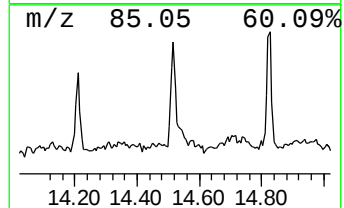
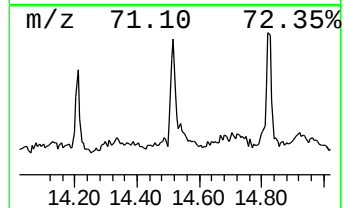
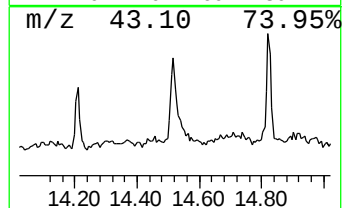
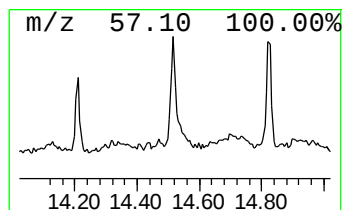
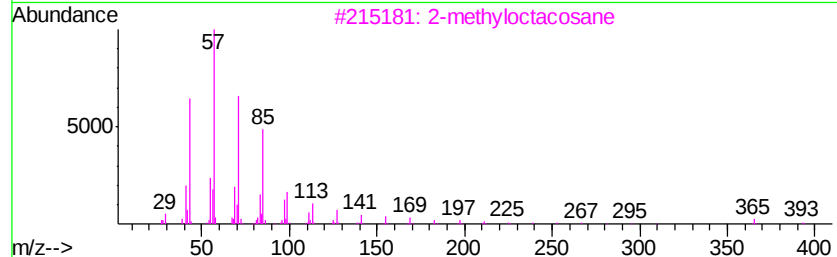
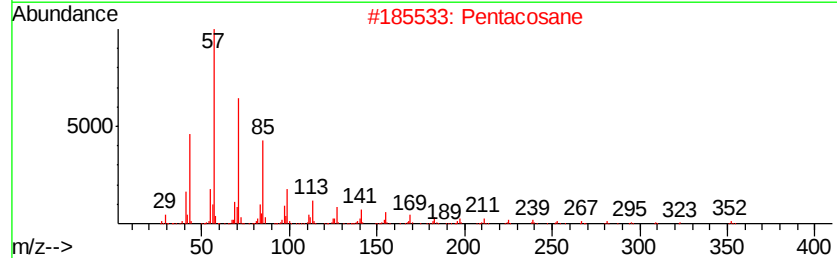
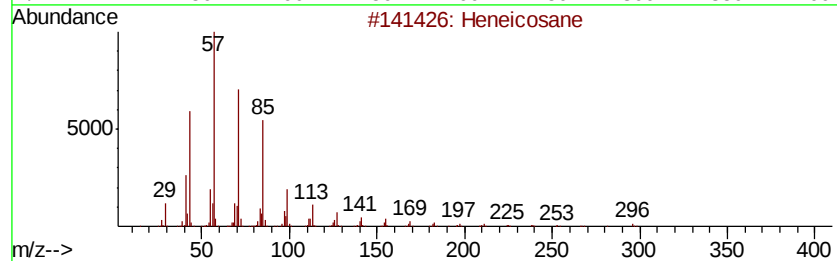
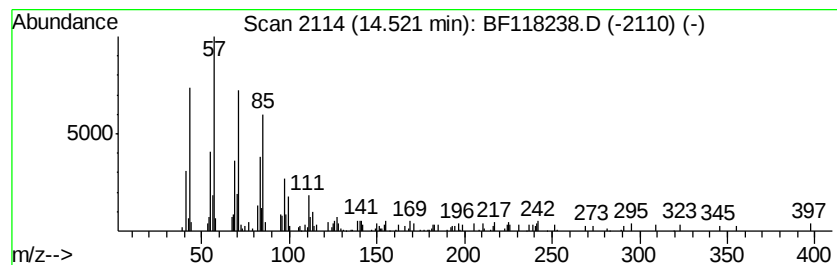
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Pentacosane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.52	5.08 ng	216112	Chrysene-d12	14.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosane	296	C21H44	000629-94-7	91
2		Pentacosane	352	C25H52	000629-99-2	91
3		2-methyloctacosane	408	C29H60	1000376-72-8	91
4		Octacosane	394	C28H58	000630-02-4	91
5		Tetratetracontane	619	C44H90	007098-22-8	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121719\
Data File : BF118238.D
Acq On : 17 Dec 2019 20:48
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ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
121619-CA-COMP

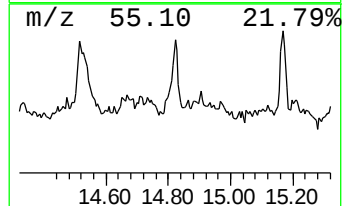
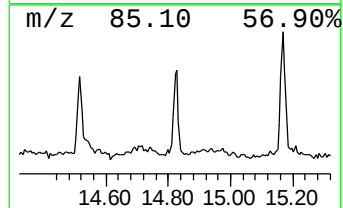
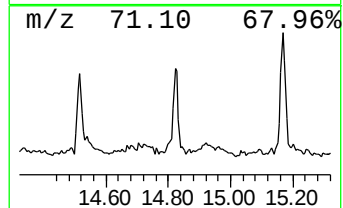
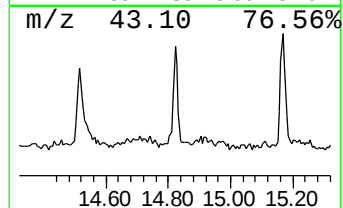
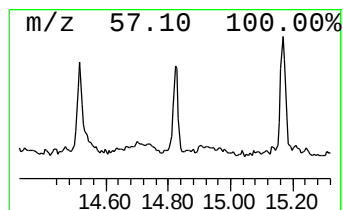
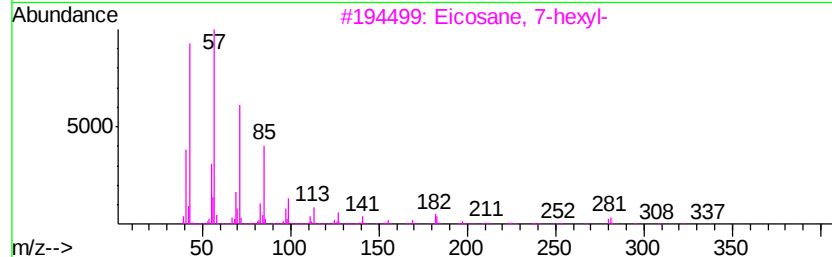
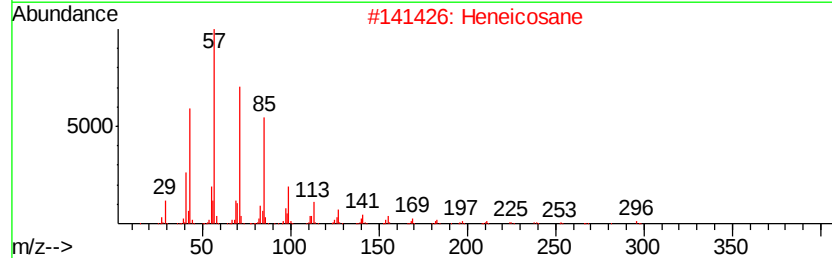
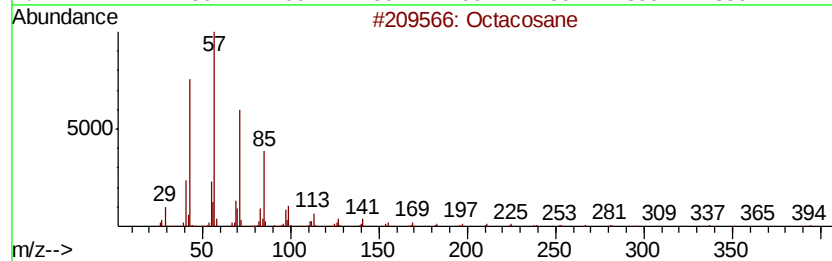
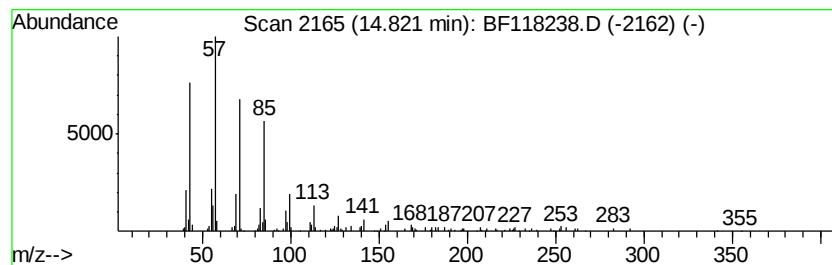
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Octacosane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.82	3.27 ng	156963	Perylene-d12	15.54

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octacosane	394	C28H58	000630-02-4	87
2		Heneicosane	296	C21H44	000629-94-7	87
3		Eicosane, 7-hexyl-	366	C26H54	055333-99-8	86
4		Eicosane, 2-methyl-	296	C21H44	001560-84-5	86
5		Octadecane	254	C18H38	000593-45-3	86



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121719\
Data File : BF118238.D
Acq On : 17 Dec 2019 20:48
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Misc :
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ClientSampleId :
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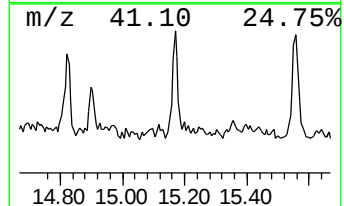
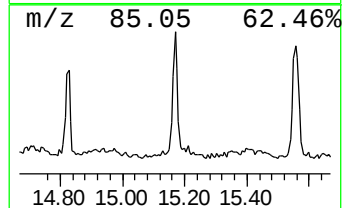
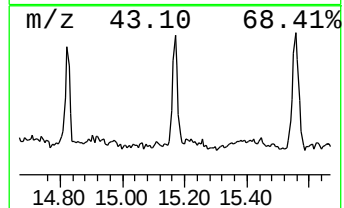
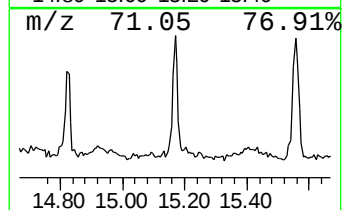
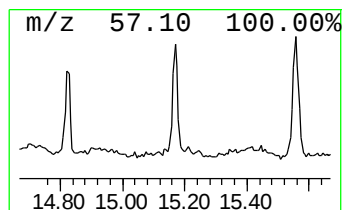
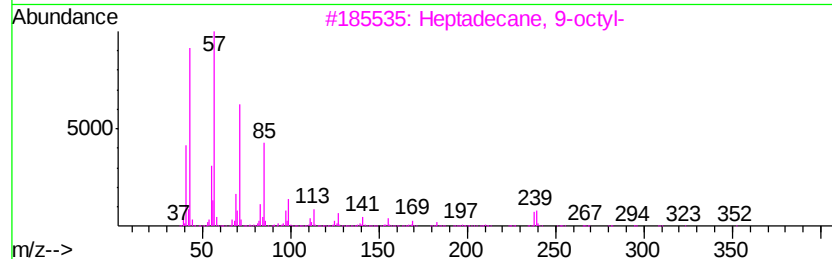
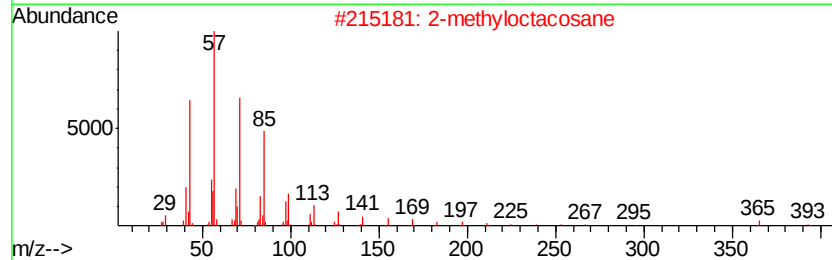
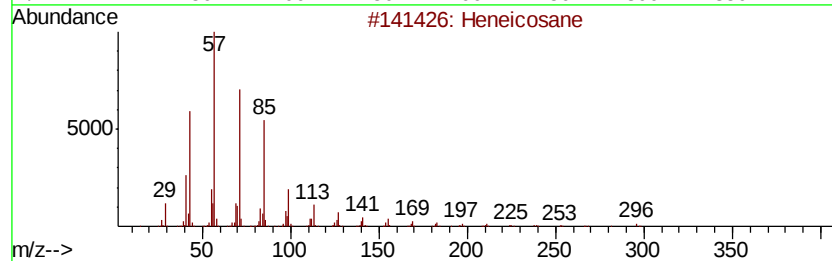
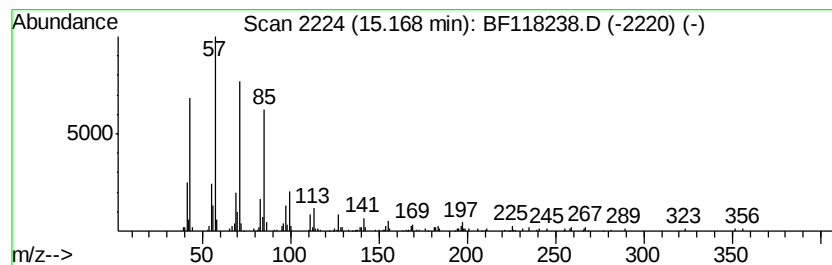
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 2-methyloctacosane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.17	4.21 ng	202270	Perylene-d12	15.54

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosane	296	C21H44	000629-94-7	91
2		2-methyloctacosane	408	C29H60	1000376-72-8	91
3		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
4		triacontane	422	C30H62	000638-68-6	91
5		Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121719\
Data File : BF118238.D
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Operator : JU
Sample : K6329-01
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
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ClientSampleId :
121619-CA-COMP

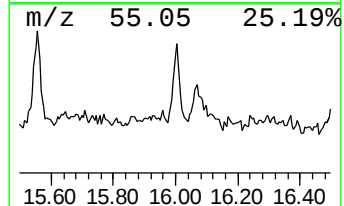
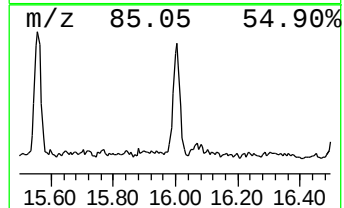
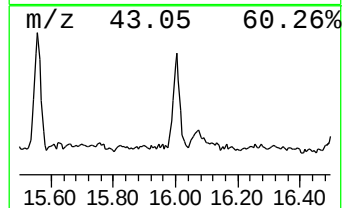
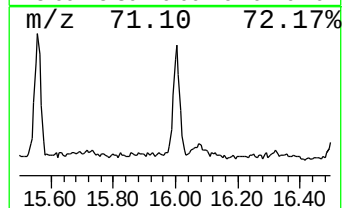
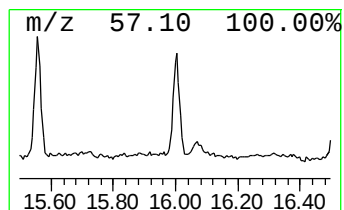
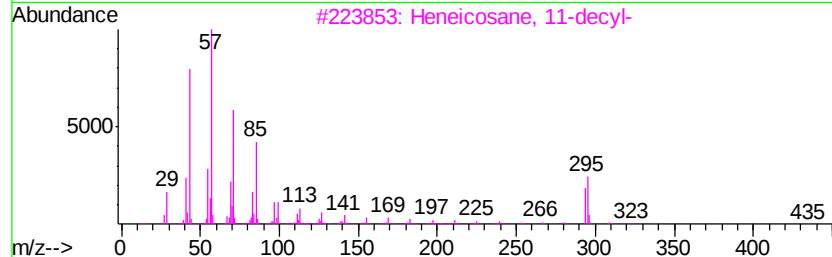
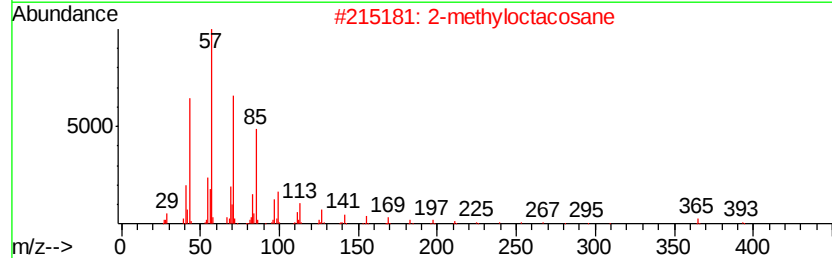
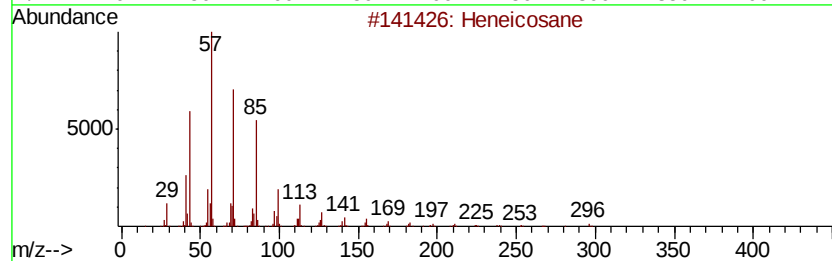
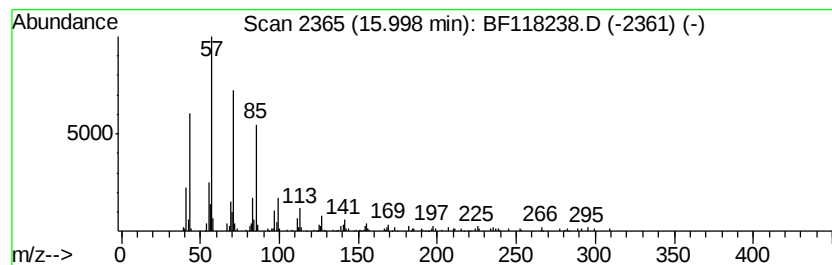
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Heneicosane, 11-decyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.00	5.47 ng	262838	Perylene-d12	15.54

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosane	296	C21H44	000629-94-7	91
2		2-methyloctacosane	408	C29H60	1000376-72-8	91
3		Heneicosane, 11-decyl-	437	C31H64	055320-06-4	91
4		Tetracontane	563	C40H82	004181-95-7	91
5		Docosane, 7-hexyl-	394	C28H58	055373-86-9	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121719\
Data File : BF118238.D
Acq On : 17 Dec 2019 20:48
Operator : JU
Sample : K6329-01
Misc :
ALS Vial : 15 Sample Multiplier: 1

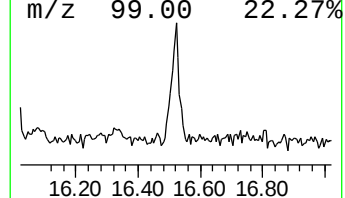
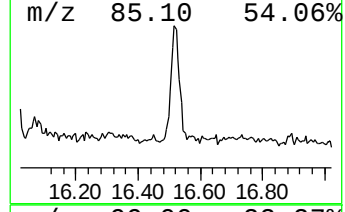
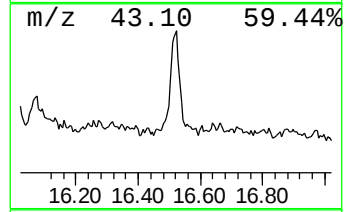
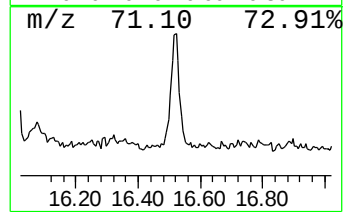
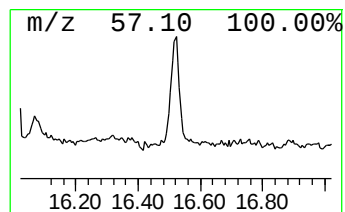
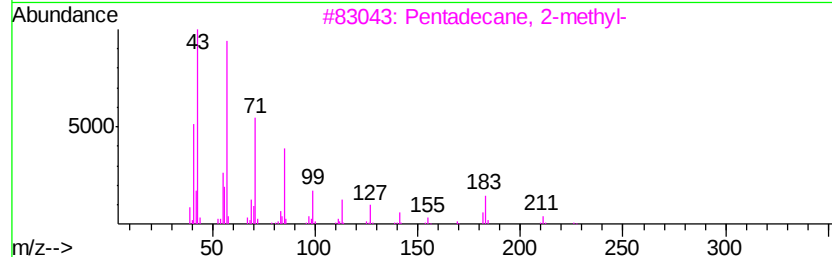
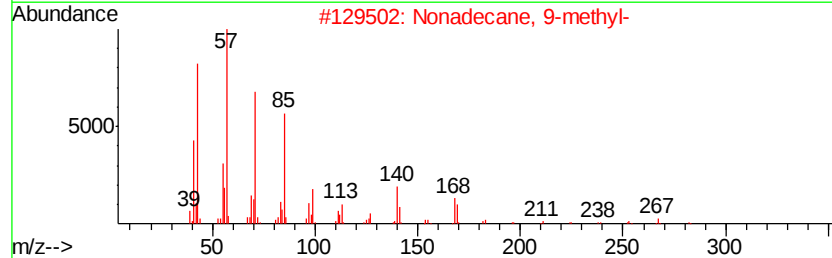
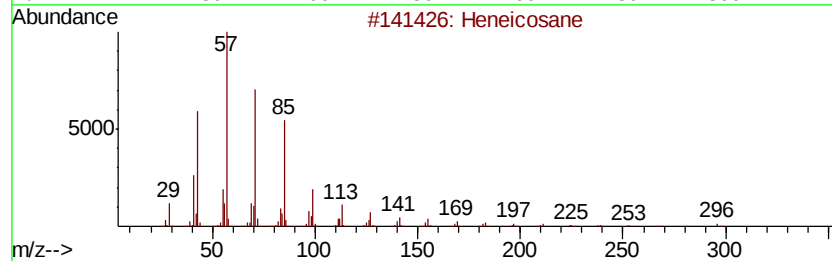
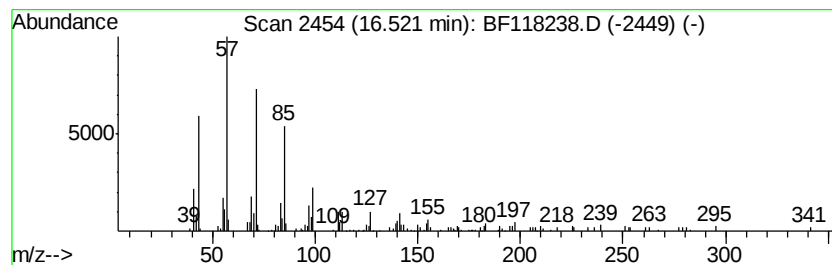
Instrument :
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ClientSampleId :
121619-CA-COMP

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 16 Nonadecane, 9-methyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD		R.T.
16.52	3.56 ng	170900	Perylene-d12		15.54
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane	296	C21H44	000629-94-7	91
2	Nonadecane, 9-methyl-	282	C20H42	013287-24-6	91
3	Pentadecane, 2-methyl-	226	C16H34	001560-93-6	87
4	Tetratetracontane	619	C44H90	007098-22-8	87
5	Tetracosane	338	C24H50	000646-31-1	86



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121719\
Data File : BF118238.D
Acq On : 17 Dec 2019 20:48
Operator : JU
Sample : K6329-01
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
121619-CA-COMP

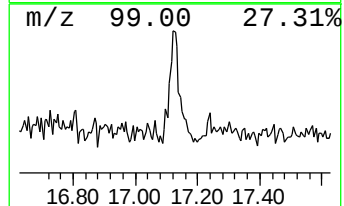
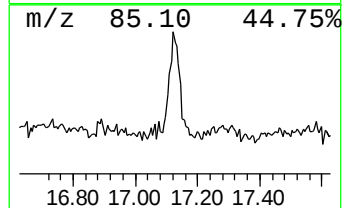
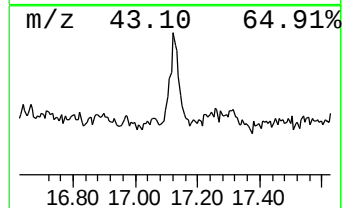
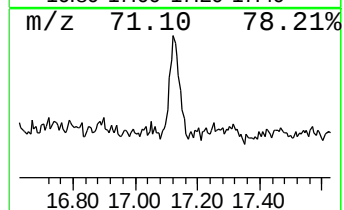
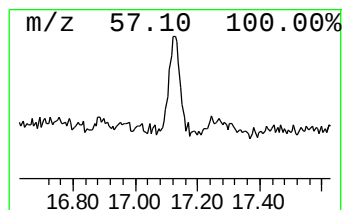
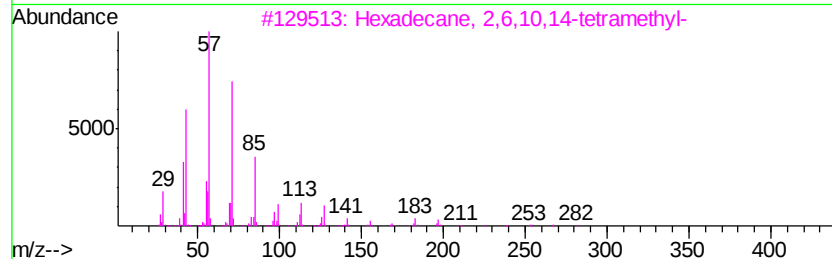
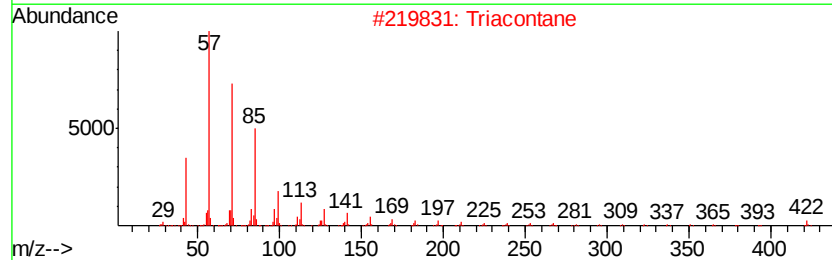
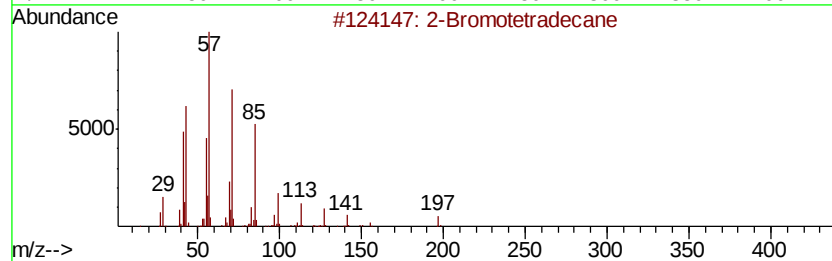
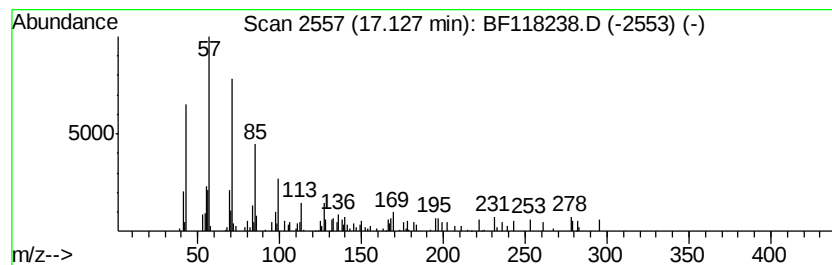
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF120619.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 2-Bromotetradecane Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.13	2.99 ng	143820	Perylene-d12	15.54

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Bromotetradecane	276	C14H29Br	074036-95-6	64
2		Triacontane	422	C30H62	000638-68-6	64
3		Hexadecane, 2,6,10,14-tetramethyl-	282	C20H42	000638-36-8	64
4		Hexacosane	366	C26H54	000630-01-3	64
5		Pentacosane	352	C25H52	000629-99-2	59



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF121719\
 Data File : BF118238.D
 Acq On : 17 Dec 2019 20:48
 Operator : JU
 Sample : K6329-01
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 121619-CA-COMP

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...	2.12	9.6	ng	373528	1	6.86	778400	20.0
2-Acetoxyisobutyr...	5.07	13.5	ng	524852	1	6.86	778400	20.0
Octane, 3,6-dimet...	6.10	3.8	ng	147823	1	6.86	778400	20.0
1R,2c,3t,4t-Tetra...	6.39	2.3	ng	87709	1	6.86	778400	20.0
unknown6.63	6.63	74.6	ng	2901810	1	6.86	778400	20.0
n-Hexadecanoic acid	11.92	5.9	ng	268120	4	11.40	911155	20.0
Pentafluoropropio...	13.89	13.5	ng	574303	5	14.06	850483	20.0
Heneicosane	14.21	2.5	ng	105284	5	14.06	850483	20.0
Pentacosane	14.52	5.1	ng	216112	5	14.06	850483	20.0
Octacosane	14.82	3.3	ng	156963	6	15.54	961109	20.0
2-methyloctacosane	15.17	4.2	ng	202270	6	15.54	961109	20.0
Heneicosane, 11-d...	16.00	5.5	ng	262838	6	15.54	961109	20.0
Nonadecane, 9-met...	16.52	3.6	ng	170900	6	15.54	961109	20.0
2-Bromotetradecane	17.13	3.0	ng	143820	6	15.54	961109	20.0