

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF060520\
 Data File : BF120561.D
 Acq On : 5 Jun 2020 15:58
 Operator : JU/CG
 Sample : L2890-03
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 72-11979-72-11969

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	1.946	7	10	21	rVV	49137	72662	2.25%	0.239%
2	2.040	21	26	34	rVB	822249	1056897	32.75%	3.475%
3	5.434	599	603	606	rBV	3320229	3226871	100.00%	10.609%
4	6.457	772	777	780	rBV	2114728	2095446	64.94%	6.889%
5	6.816	834	838	841	rBV	952977	732694	22.71%	2.409%
6	6.975	861	865	868	rBV	2324569	1941096	60.15%	6.382%
7	7.381	928	934	937	rBV	1721693	1489505	46.16%	4.897%
8	8.098	1052	1056	1059	rBV	1232189	952972	29.53%	3.133%
9	9.122	1227	1230	1233	rVB	73144	59054	1.83%	0.194%
10	9.175	1234	1239	1242	rBV	3344371	2781146	86.19%	9.144%
11	9.257	1249	1253	1257	rBV4	50636	90658	2.81%	0.298%
12	9.569	1302	1306	1310	rBV2	560234	640802	19.86%	2.107%
13	9.669	1321	1323	1328	rVB2	52912	58397	1.81%	0.192%
14	9.722	1328	1332	1334	rBV3	118201	142446	4.41%	0.468%
15	9.763	1336	1339	1342	rVB	178528	186995	5.79%	0.615%
16	9.851	1348	1354	1358	rBV	1195172	1161193	35.99%	3.818%
17	9.892	1358	1361	1363	rBV3	80363	91275	2.83%	0.300%
18	10.045	1385	1387	1391	rBV2	84505	94590	2.93%	0.311%
19	10.110	1395	1398	1401	rBV2	230659	241129	7.47%	0.793%
20	10.169	1406	1408	1409	rBV	85889	69577	2.16%	0.229%
21	10.228	1415	1418	1421	rBV2	124272	147177	4.56%	0.484%
22	10.263	1421	1424	1428	rBV3	228144	269270	8.34%	0.885%
23	10.510	1463	1466	1470	rVB	947394	886569	27.47%	2.915%
24	10.645	1485	1489	1491	rBV	1347275	1219525	37.79%	4.009%
25	10.780	1507	1512	1515	rBV	2255649	2491180	77.20%	8.190%
26	11.245	1587	1591	1597	rVB	2337600	2408711	74.65%	7.919%
27	11.345	1605	1608	1618	rBV2	1011176	1204771	37.34%	3.961%
28	11.592	1647	1650	1653	rBV	600024	564461	17.49%	1.856%
29	12.322	1772	1774	1779	rVB	316915	362068	11.22%	1.190%
30	12.780	1850	1852	1855	rVB	209618	158457	4.91%	0.521%
31	12.927	1873	1877	1880	rBV	2317628	2104096	65.21%	6.918%
32	13.980	2052	2056	2058	rBV2	740221	748192	23.19%	2.460%
33	15.427	2298	2302	2305	rBV	557205	666267	20.65%	2.191%

LSC Area Percent Report

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Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
72-11979-72-11969

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-BF052820.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

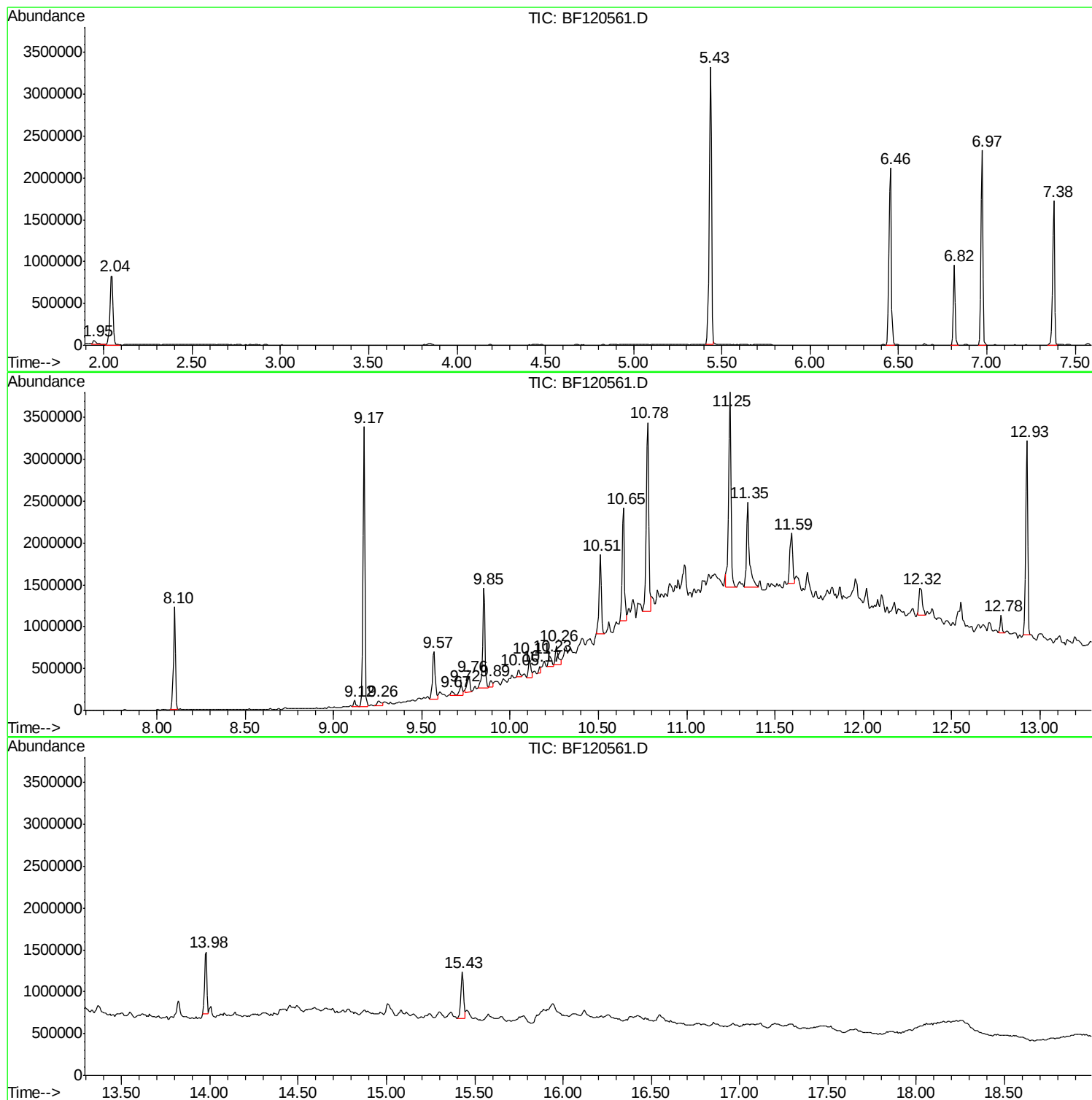
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF052820.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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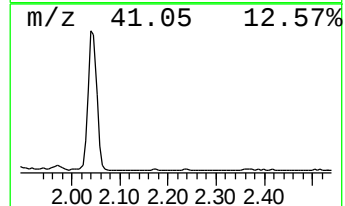
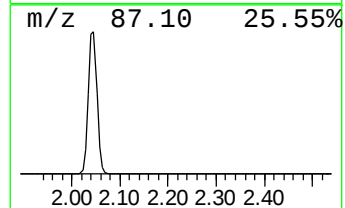
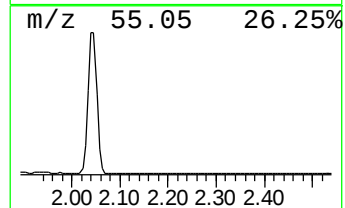
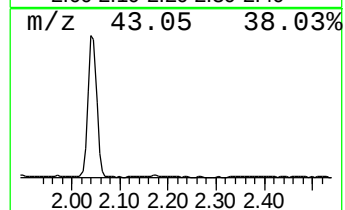
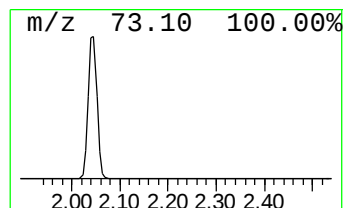
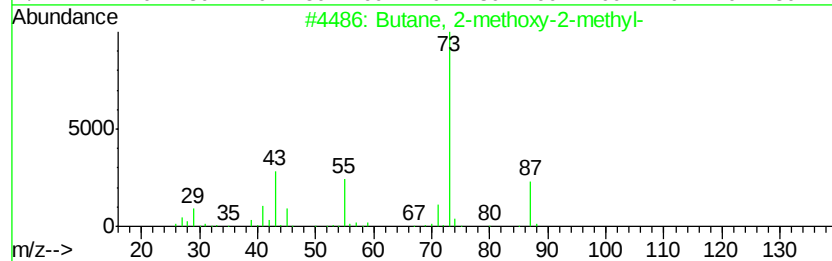
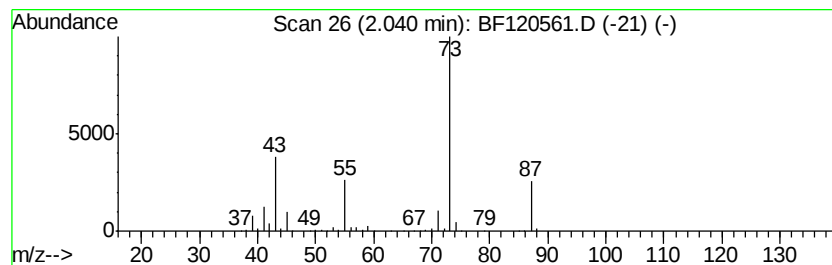
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.04	28.85 ng	1056900	1,4-Dichlorobenzene-d4	6.82

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	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		1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	37
4		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	17
5		Silane, tetramethyl-	88	C4H12Si	000075-76-3	12



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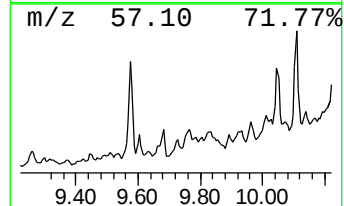
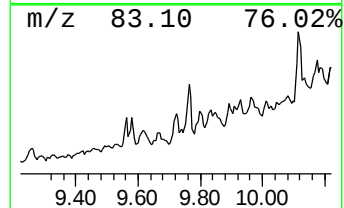
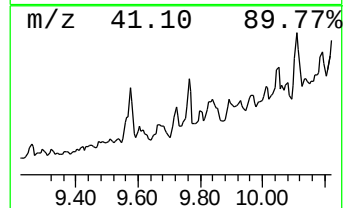
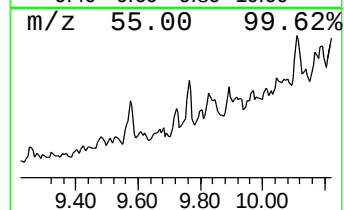
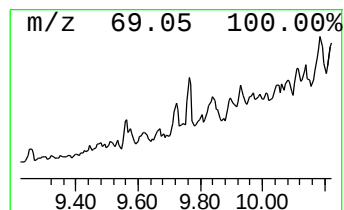
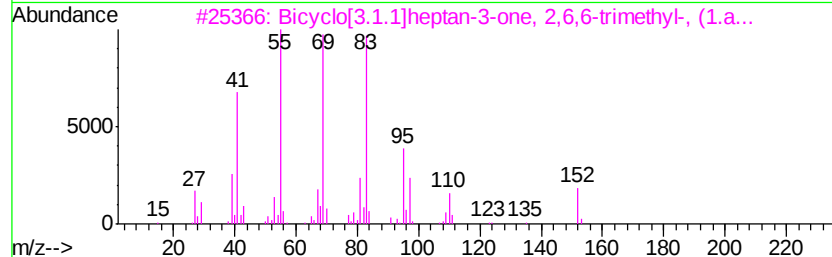
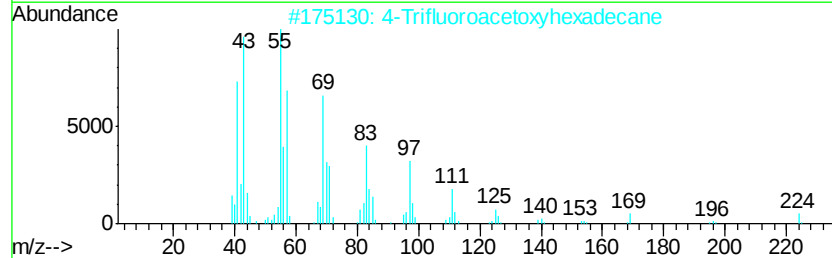
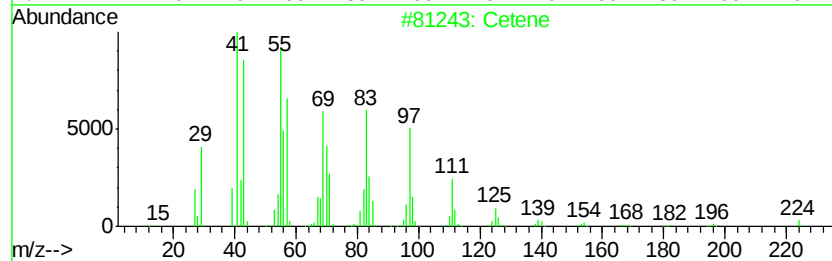
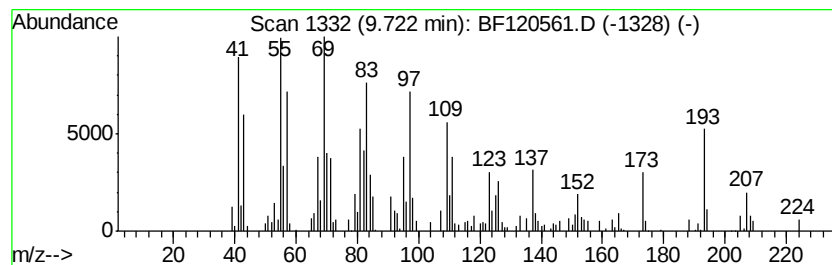
Instrument :
BNA_F
ClientSampleId :
72-11979-72-11969

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 3 Cetene Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.72	2.45 ng	142446	Acenaphthene-d10	9.85
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Cetene	224 C16H32	000629-73-2 55
2		4-Trifluoroacetoxyhexadecane	338 C18H33F3O2	1000215-97-5 53
3		Bicyclo[3.1.1]heptan-3-one, 2,6,...	152 C10H16O	015358-88-0 50
4		3-Hexadecene, (Z)-	224 C16H32	034303-81-6 50
5		Cyclohexadecane	224 C16H32	000295-65-8 50



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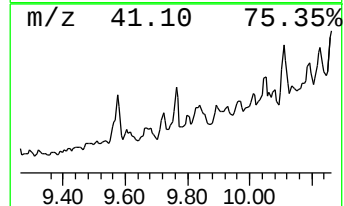
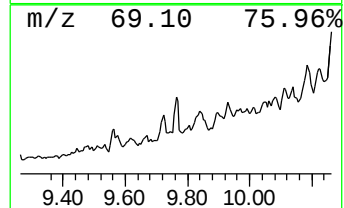
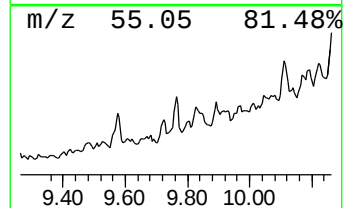
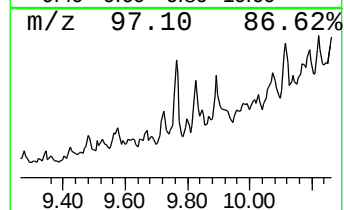
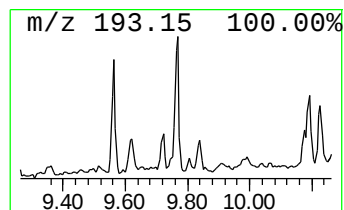
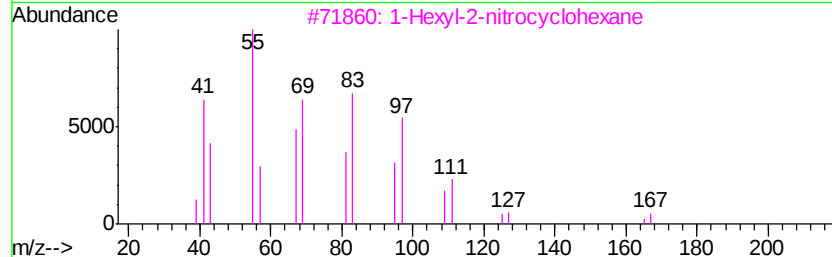
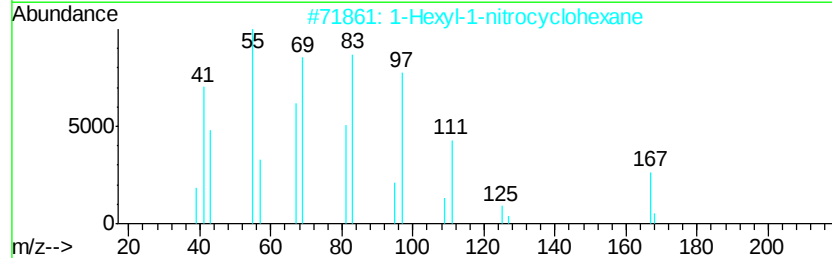
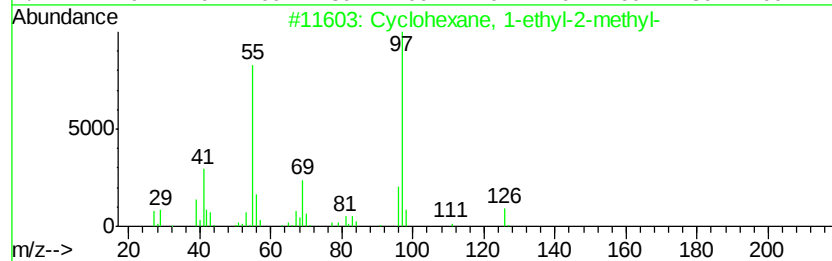
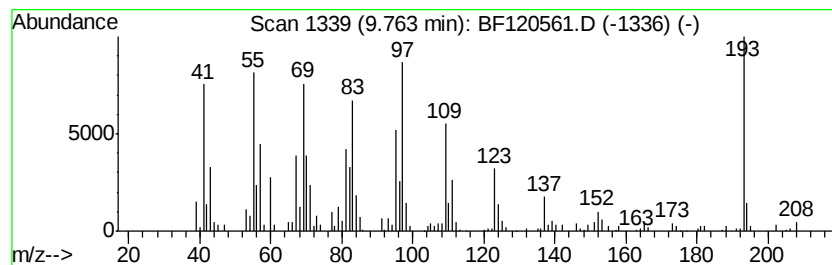
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 unknown9.76 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.76	3.22 ng	186995	Acenaphthene-d10	9.85

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexane, 1-ethyl-2-methyl-	126	C9H18	003728-54-9	35
2		1-Hexyl-1-nitrocyclohexane	213	C12H23NO2	118252-09-8	35
3		1-Hexyl-2-nitrocyclohexane	213	C12H23NO2	118252-04-3	35
4		Hexahydro-1-oxa-cyclopropa[d]ind...	152	C9H12O2	037643-23-5	30
5		Cyclopentane, (1-methylbutyl)-	140	C10H20	004737-43-3	25



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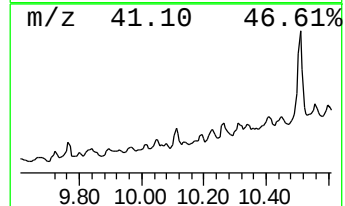
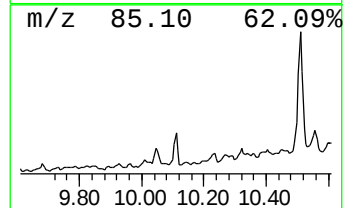
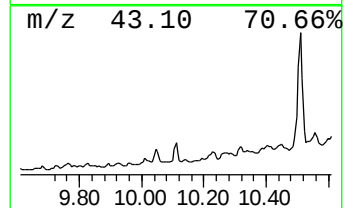
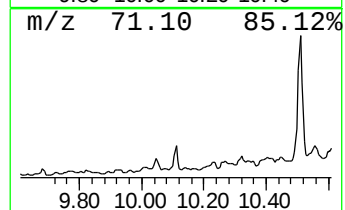
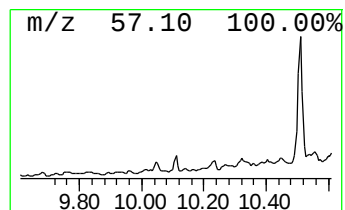
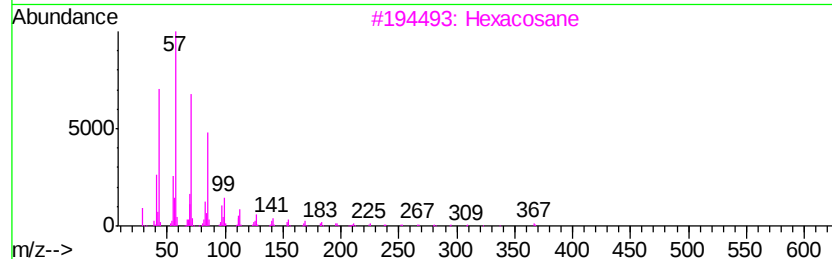
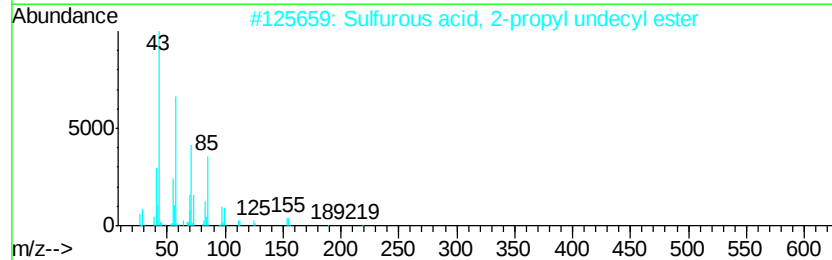
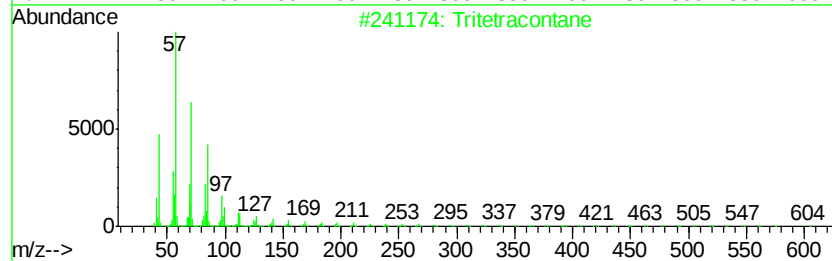
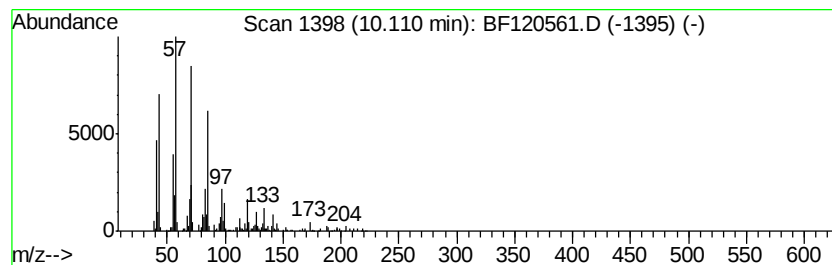
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Tritetracontane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.11	4.15 ng	241129	Acenaphthene-d10	9.85	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tritetracontane	605	C43H88	007098-21-7	80
2	Sulfurous acid, 2-propyl undecyl...	278	C14H30O3S	1000309-12-2	72
3	Hexacosane	366	C26H54	000630-01-3	72
4	Tetratetracontane	619	C44H90	007098-22-8	64
5	Nonadecane	268	C19H40	000629-92-5	58



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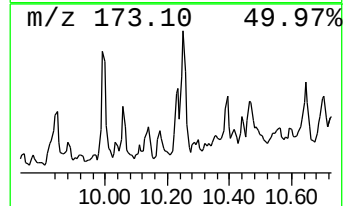
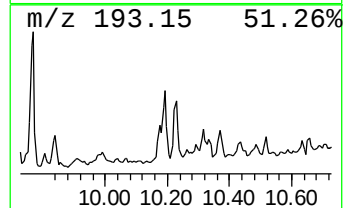
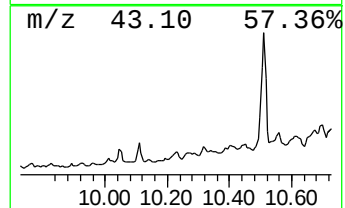
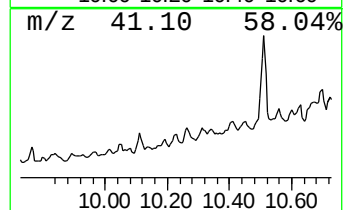
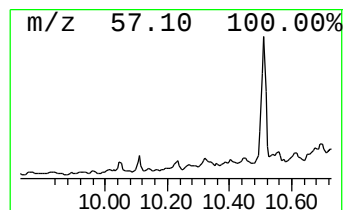
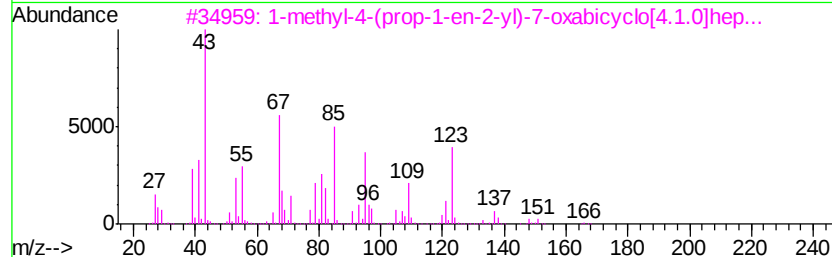
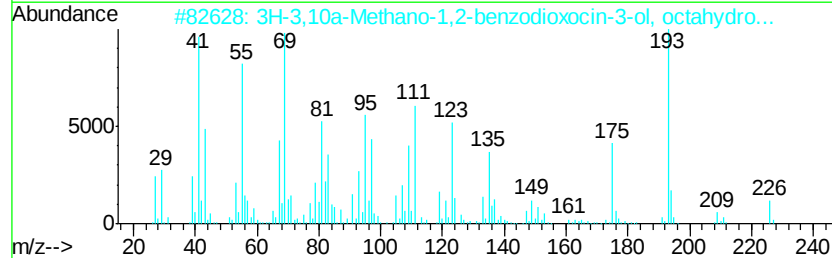
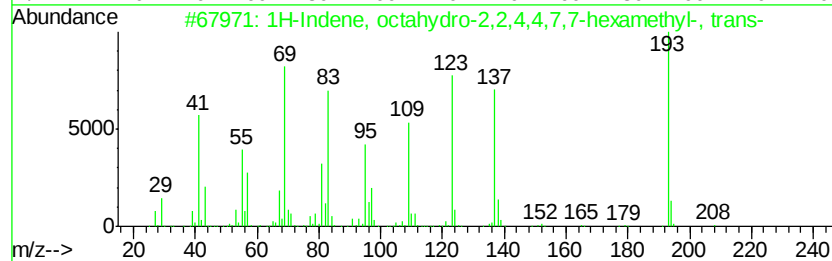
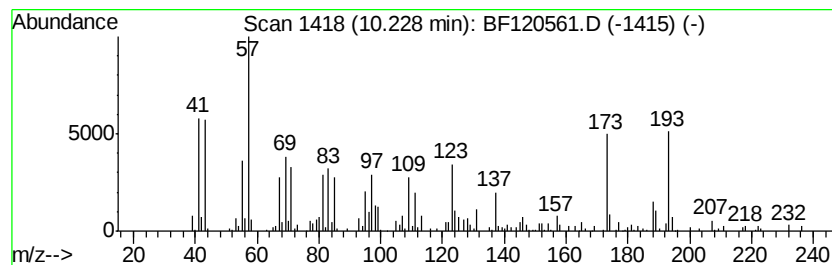
Instrument :
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 unknown10.23 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.23	2.53 ng	147177	Acenaphthene-d10	9.85	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, octahydro-2,2,4,4,7,7...	208	C15H28	054832-83-6	38
2	3H-3,10a-Methano-1,2-benzodioxoc...	226	C13H22O3	095906-83-5	27
3	1-methyl-4-(prop-1-en-2-yl)-7-ox...	166	C10H14O2	1000365-66-7	14
4	2-Piperidinone, N-[4-bromo-n-but...	233	C9H16BrNO	195194-80-0	14
5	1-(p-Fluorophenyl)-4-piperidone	193	C11H12FNO	1000238-56-7	14



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ALS Vial : 10 Sample Multiplier: 1

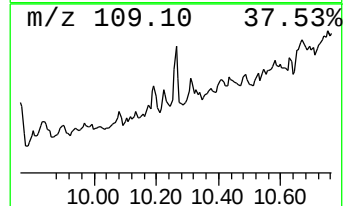
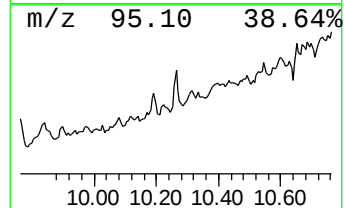
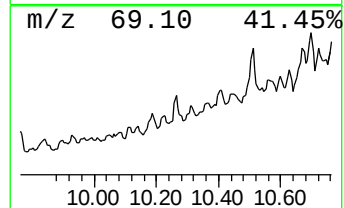
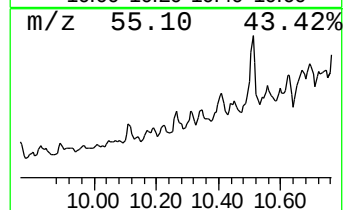
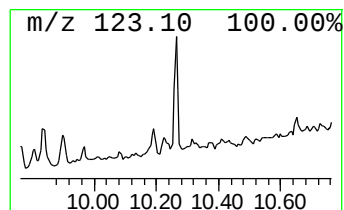
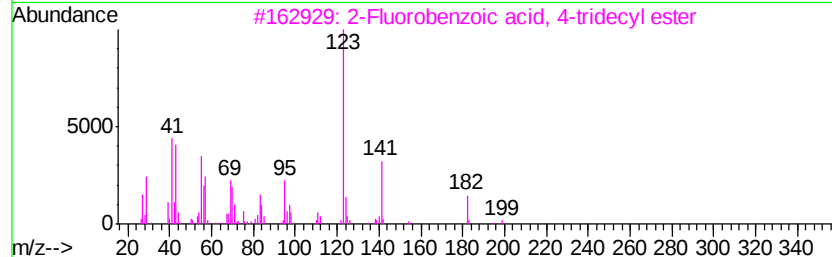
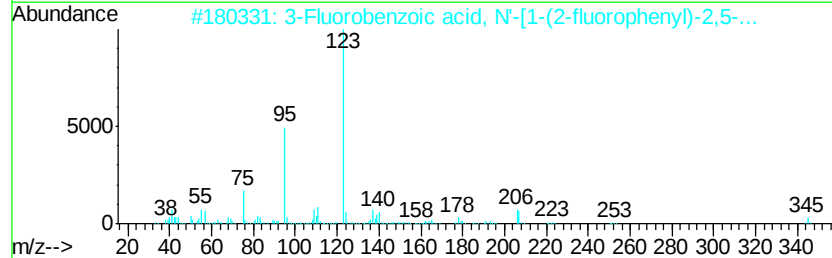
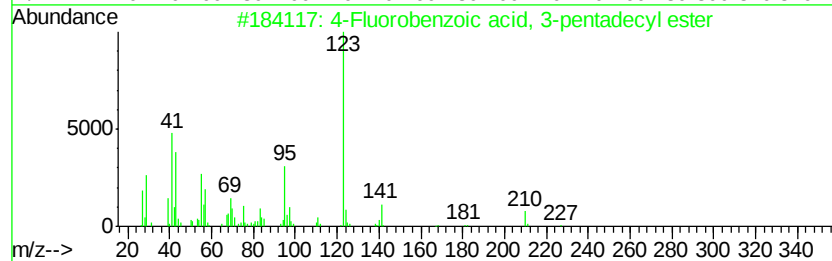
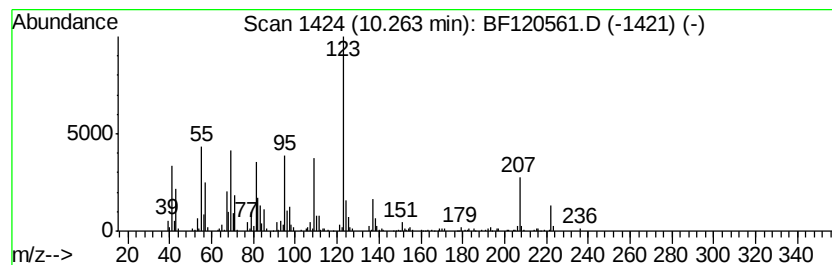
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 4-Fluorobenzoic acid, 3-pen... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD		R.T.	
10.26	4.64 ng	269270	Acenaphthene-d10		9.85	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		4-Fluorobenzoic acid, 3-pentadec...	350	C22H35F02	1000280-68-4	53
2		3-Fluorobenzoic acid, N'-[1-(2-f...	345	C17H13F2N3O3	1000304-52-3	50
3		2-Fluorobenzoic acid, 4-tridecyl...	322	C20H31F02	1000280-61-2	47
4		N-(4,6-Dimethyl-pyrimidin-2-yl)-...	393	C21H20FN5O2	1000300-88-4	47
5		1,3-Diazocin-2-one, perhydro-1-(...	250	C13H15FN2O2	1000267-45-2	47



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF060520\
Data File : BF120561.D
Acq On : 5 Jun 2020 15:58
Operator : JU/CG
Sample : L2890-03
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
72-11979-72-11969

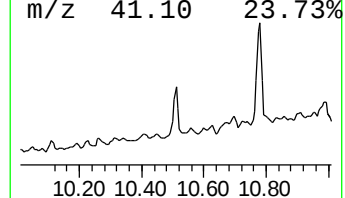
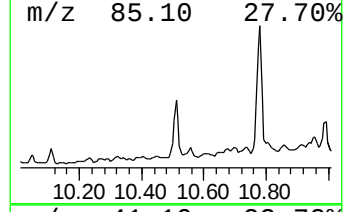
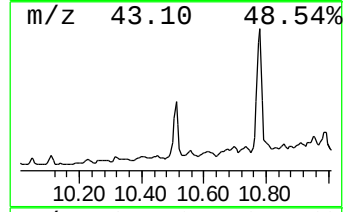
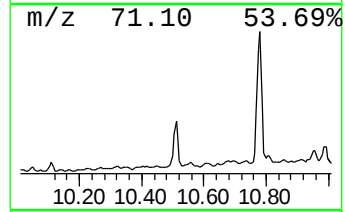
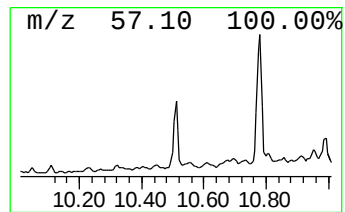
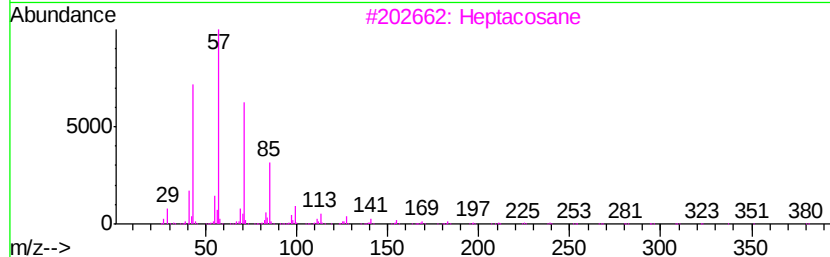
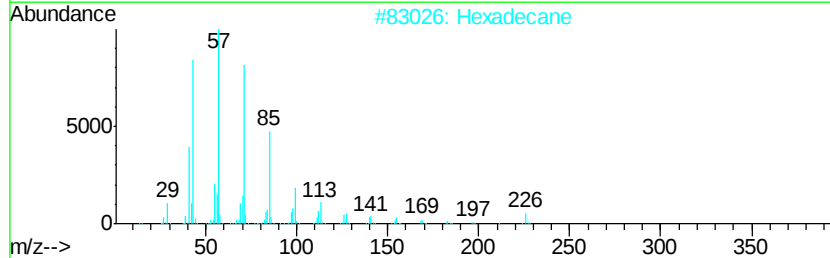
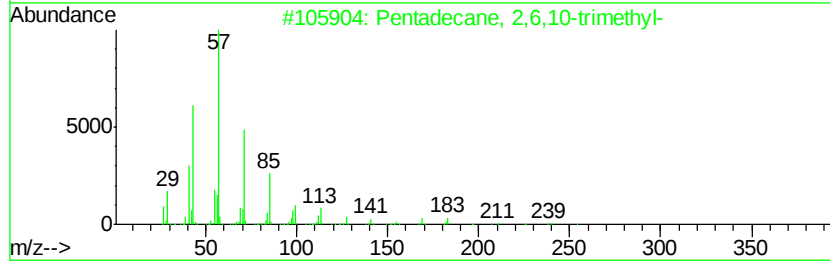
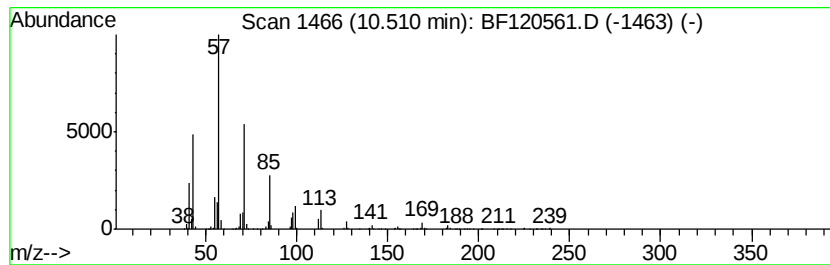
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Pentadecane, 2,6,10-trimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.51	15.27 ng	886569	Acenaphthene-d10	9.85

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentadecane, 2,6,10-trimethyl-	254	C18H38	003892-00-0	90
2		Hexadecane	226	C16H34	000544-76-3	87
3		Heptacosane	380	C27H56	000593-49-7	86
4		Undecane, 3,5-dimethyl-	184	C13H28	017312-81-1	81
5		Heneicosane	296	C21H44	000629-94-7	80



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF060520\
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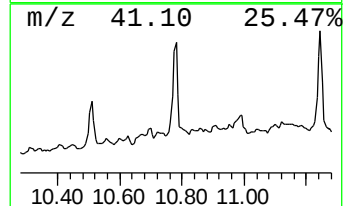
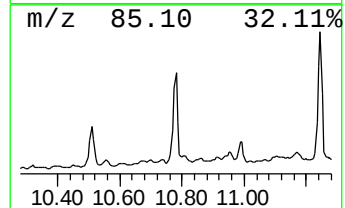
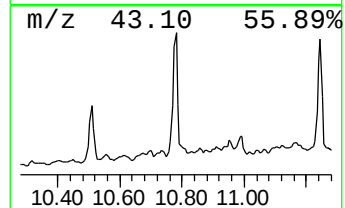
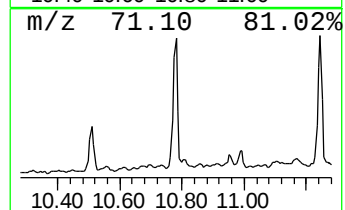
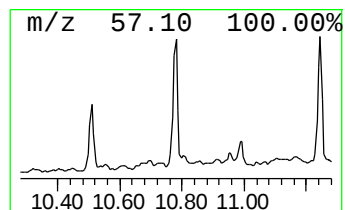
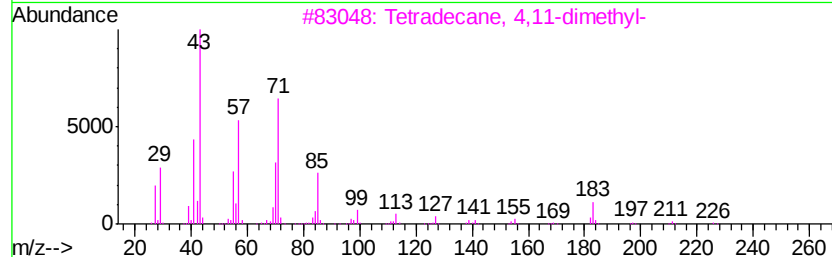
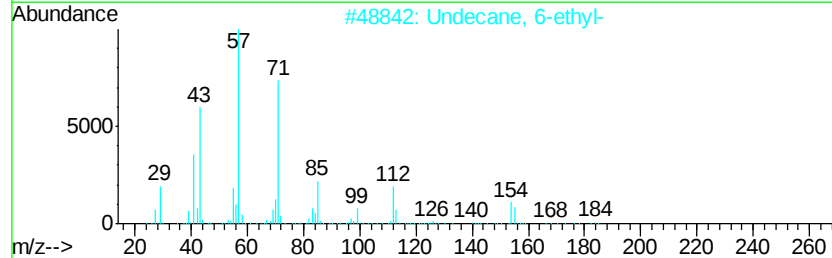
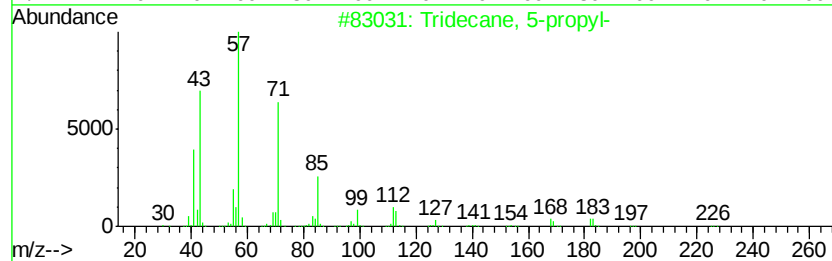
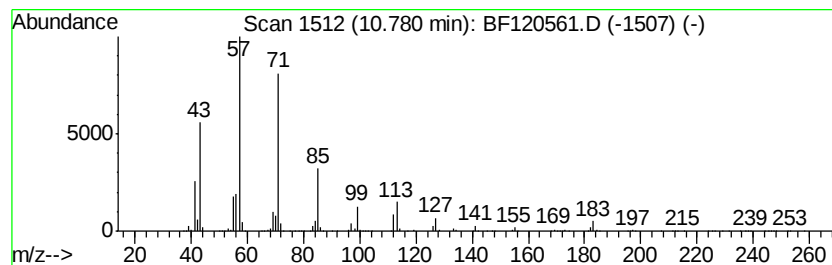
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF052820.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Tridecane, 5-propyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.78	41.36 ng	2491180	Phenanthrene-d10	11.35
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Tridecane, 5-propyl-	226 C16H34	055045-11-9 87
2		Undecane, 6-ethyl-	184 C13H28	017312-60-6 80
3		Tetradecane, 4,11-dimethyl-	226 C16H34	055045-12-0 72
4		Sulfurous acid, decyl 2-ethylhex...	334 C18H38O3S	1000309-19-3 72
5		Sulfurous acid, 2-ethylhexyl non...	320 C17H36O3S	1000309-19-2 72



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF060520\
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Sample : L2890-03
Misc :
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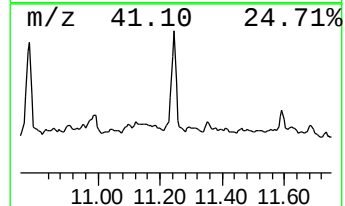
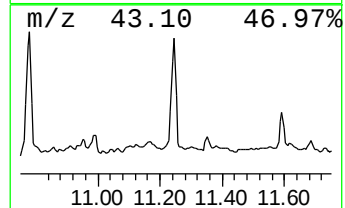
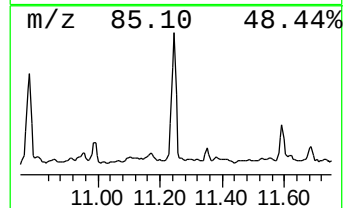
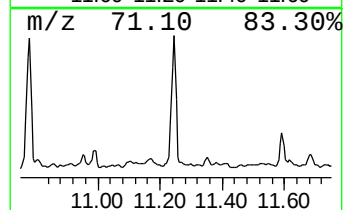
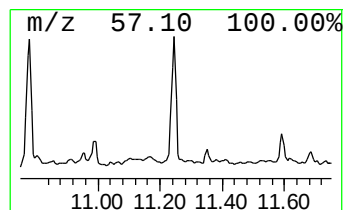
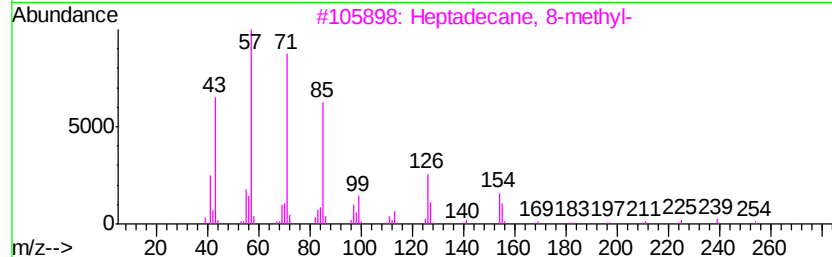
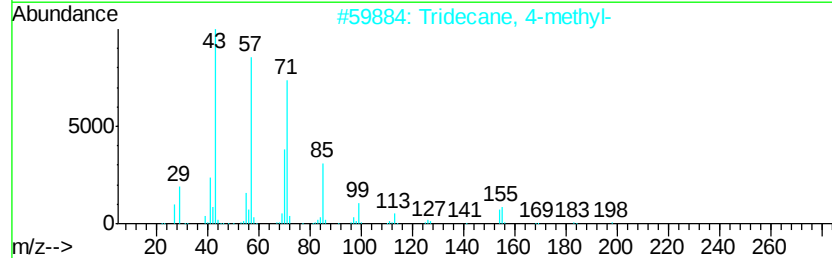
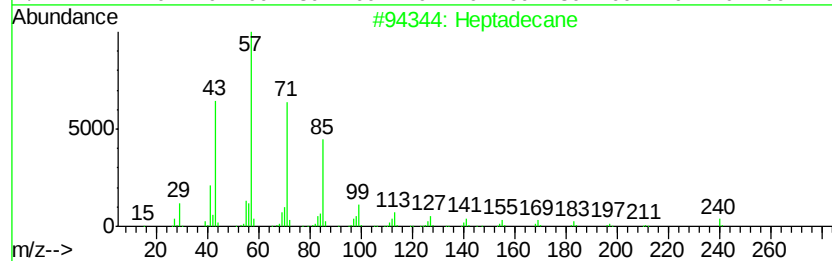
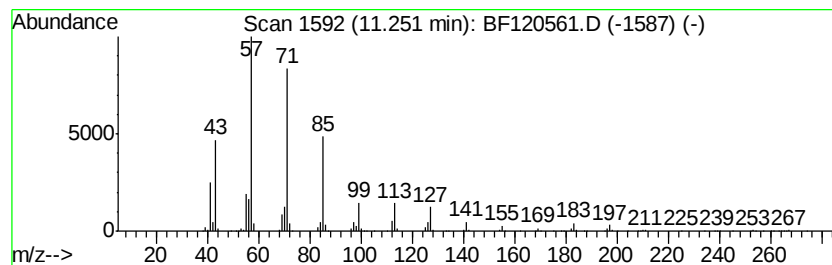
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Heptadecane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.25	39.99 ng	2408710	Phenanthrene-d10	11.35
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Heptadecane	240 C17H36	000629-78-7	90
2	Tridecane, 4-methyl-	198 C14H30	026730-12-1	87
3	Heptadecane, 8-methyl-	254 C18H38	013287-23-5	83
4	Dodecane, 4-methyl-	184 C13H28	006117-97-1	81
5	2-Bromotetradecane	276 C14H29Br	074036-95-6	80



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF060520\
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Sample : L2890-03
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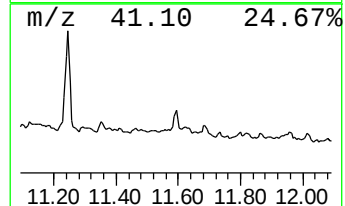
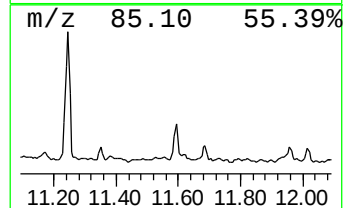
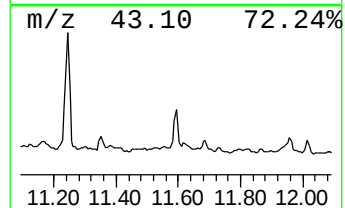
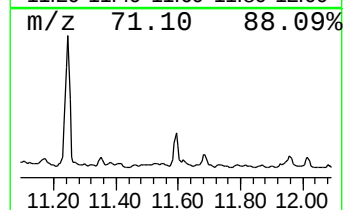
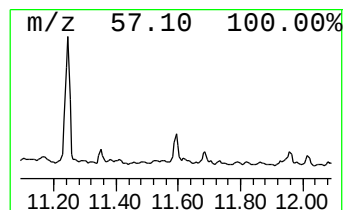
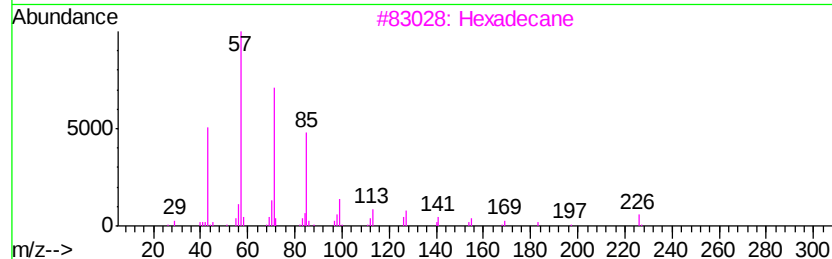
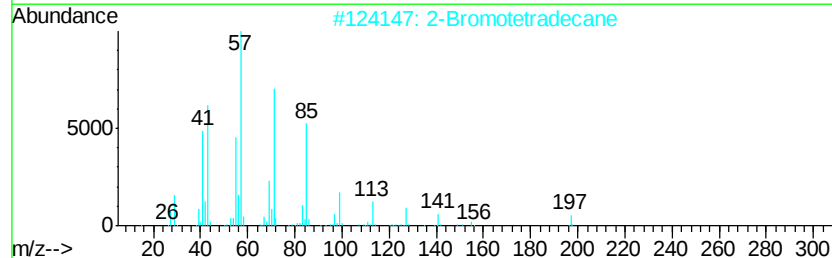
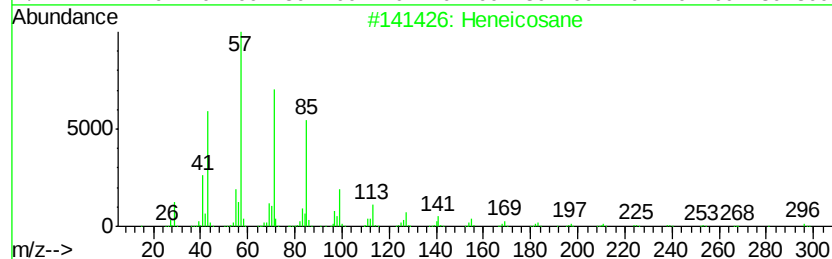
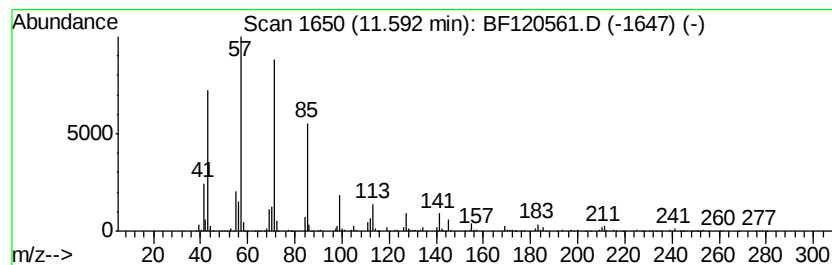
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Heneicosane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD		R.T.
11.59	9.37 ng	564461	Phenanthrene-d10		11.35
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane	296	C21H44	000629-94-7	91
2	2-Bromotetradecane	276	C14H29Br	074036-95-6	90
3	Hexadecane	226	C16H34	000544-76-3	90
4	Tetracosane	338	C24H50	000646-31-1	86
5	Octadecane	254	C18H38	000593-45-3	86



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF060520\
Data File : BF120561.D
Acq On : 5 Jun 2020 15:58
Operator : JU/CG
Sample : L2890-03
Misc :
ALS Vial : 10 Sample Multiplier: 1

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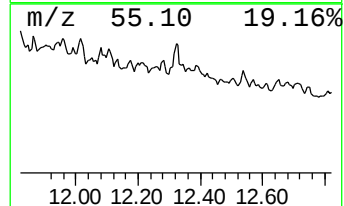
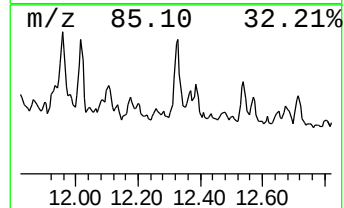
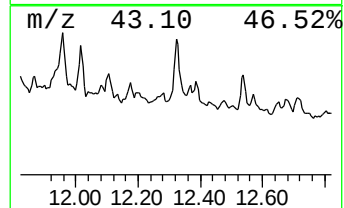
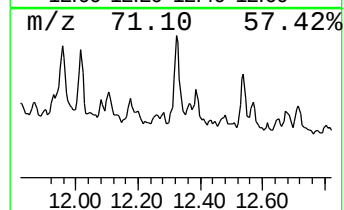
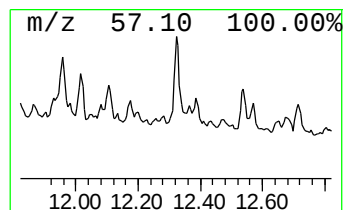
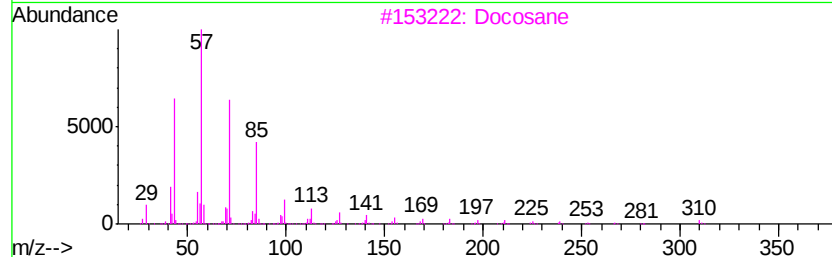
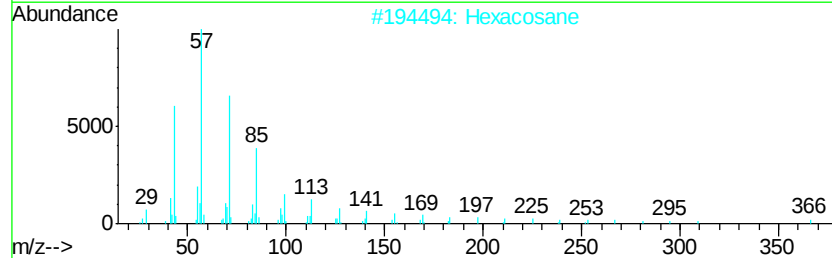
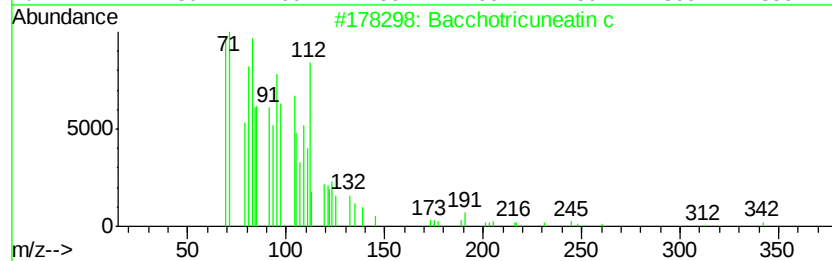
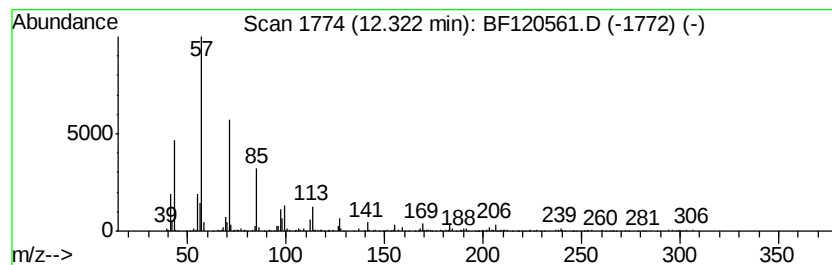
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Bacchotricuneatin c Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.32	6.01 ng	362068	Phenanthrene-d10	11.35

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Bacchotricuneatin c	342	C20H22O5	066563-30-2	97
2		Hexacosane	366	C26H54	000630-01-3	90
3		Docosane	310	C22H46	000629-97-0	86
4		Pentadecane, 2,6,10-trimethyl-	254	C18H38	003892-00-0	86
5		Dodecane	170	C12H26	000112-40-3	74



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF060520\
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 Acq On : 5 Jun 2020 15:58
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 Sample : L2890-03
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 ALS Vial : 10 Sample Multiplier: 1

Instrument :
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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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
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Cetene	9.72	2.5	ng	142446	3	9.85	1161190	20.0
unknown9.76	9.76	3.2	ng	186995	3	9.85	1161190	20.0
Tritetracontane	10.11	4.2	ng	241129	3	9.85	1161190	20.0
unknown10.23	10.23	2.5	ng	147177	3	9.85	1161190	20.0
4-Fluorobenzoic a...	10.26	4.6	ng	269270	3	9.85	1161190	20.0
Pentadecane, 2,6,...	10.51	15.3	ng	886569	3	9.85	1161190	20.0
Tridecane, 5-propyl-	10.78	41.4	ng	2491180	4	11.35	1204770	20.0
Heptadecane	11.25	40.0	ng	2408710	4	11.35	1204770	20.0
Heneicosane	11.59	9.4	ng	564461	4	11.35	1204770	20.0
Bacchotricuneatin c	12.32	6.0	ng	362068	4	11.35	1204770	20.0