

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP062620\
 Data File : BP002556.D
 Acq On : 26 Jun 2020 17:56
 Operator : CG/JU
 Sample : L3114-01
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 VNJ-242

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_P\METHODS\8270-BP060520.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	5.205	387	394	410	rBV	2801624	4511107	57.82%	6.138%
2	6.481	603	611	621	rVB2	48072	98819	1.27%	0.134%
3	6.775	649	661	675	rBV	2945425	4917662	63.03%	6.691%
4	7.581	781	798	807	rBV	776010	1234174	15.82%	1.679%
5	7.899	844	852	862	rBV	2732695	4500625	57.68%	6.123%
6	8.728	985	993	1001	rBV	2059009	3421567	43.85%	4.655%
7	10.352	1261	1269	1281	rBV	993930	1725475	22.12%	2.348%
8	11.422	1445	1451	1452	rBV5	63286	105933	1.36%	0.144%
9	12.116	1565	1569	1573	rBV3	172911	307004	3.93%	0.418%
10	12.593	1646	1650	1656	rBV6	147634	302521	3.88%	0.412%
11	12.663	1657	1662	1667	rBV	593724	994856	12.75%	1.354%
12	12.793	1682	1684	1686	rBV2	121282	133450	1.71%	0.182%
13	12.845	1686	1693	1699	rVV	4277866	7340184	94.08%	9.987%
14	13.022	1718	1723	1731	rBV	1161713	1943197	24.91%	2.644%
15	13.116	1737	1739	1744	rVB2	266991	322992	4.14%	0.439%
16	13.569	1812	1816	1820	rBV2	1061544	1692578	21.69%	2.303%
17	13.669	1827	1833	1842	rBV2	4153092	7802206	100.00%	10.615%
18	13.904	1870	1873	1878	rBV3	932194	1266236	16.23%	1.723%
19	13.987	1882	1887	1894	rBV4	2057356	5031942	64.49%	6.846%
20	14.087	1899	1904	1911	rVB	4325021	7500269	96.13%	10.205%
21	14.157	1912	1916	1921	rVV3	1167156	1866178	23.92%	2.539%
22	14.216	1922	1926	1942	rVB4	2402277	6704831	85.94%	9.122%
23	14.345	1944	1948	1956	rBV	1687696	3362015	43.09%	4.574%
24	15.057	2065	2069	2076	rVB2	2637907	4449010	57.02%	6.053%
25	23.304	3466	3471	3483	rVB2	695291	1384996	17.75%	1.884%
26	23.868	3564	3567	3575	rVB	325609	579274	7.42%	0.788%

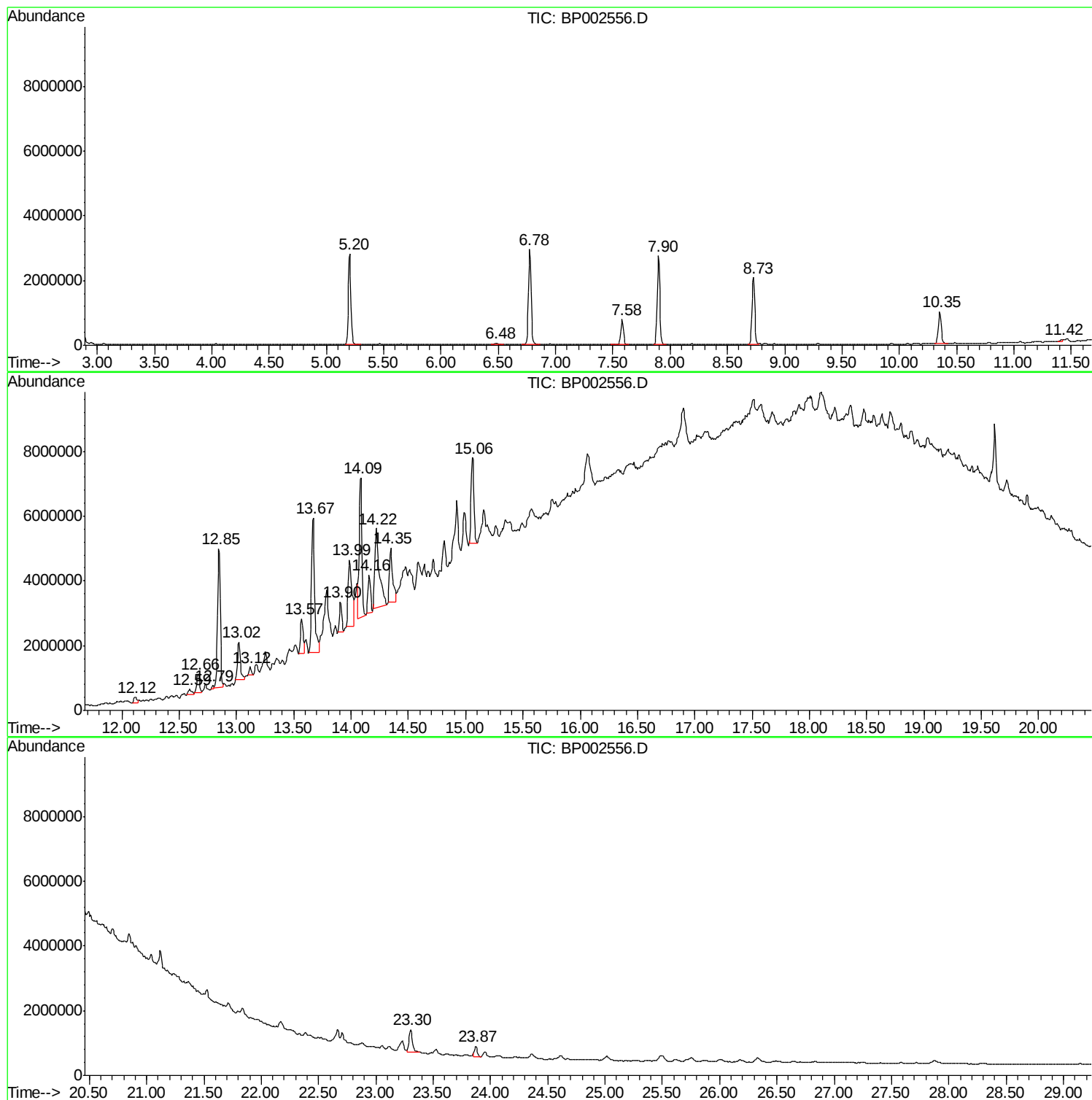
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Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP060520.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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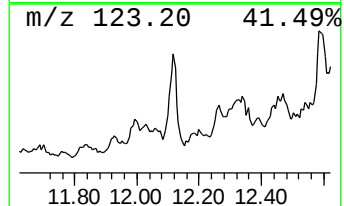
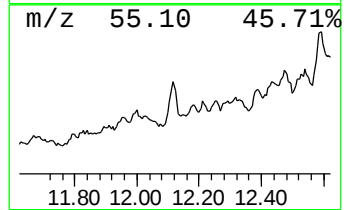
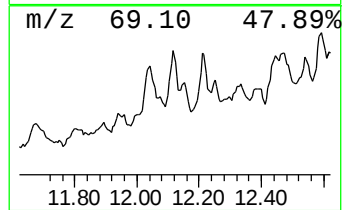
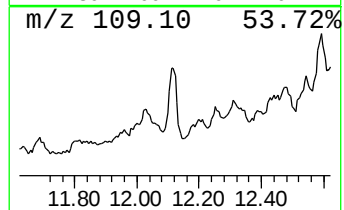
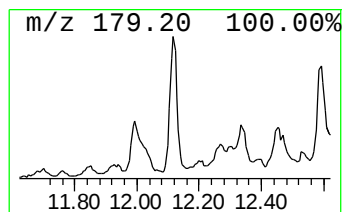
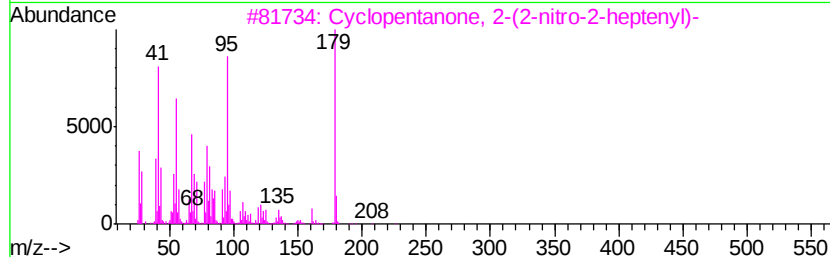
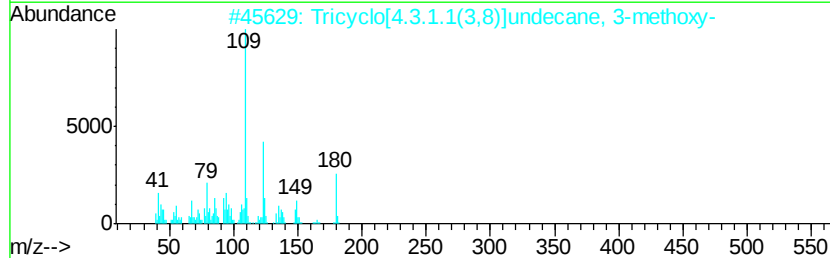
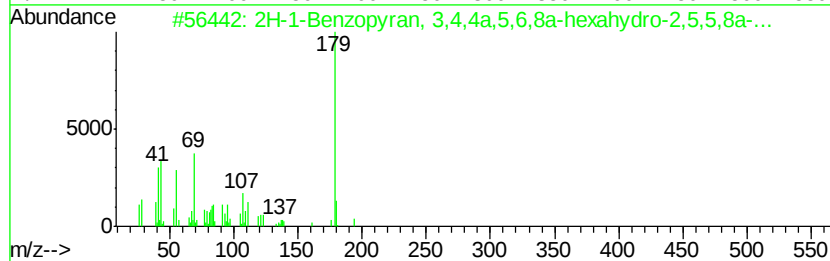
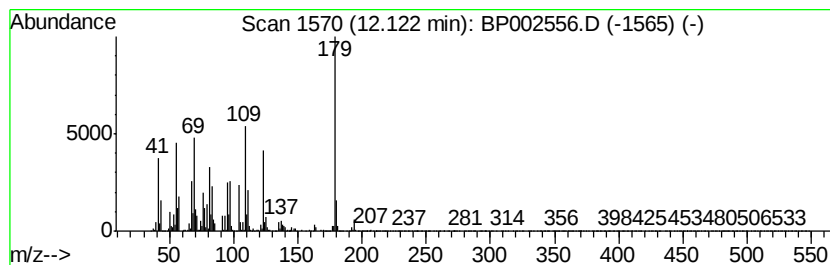
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 unknown12.12 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.12	3.56 ng	307004	Naphthalene-d8	10.35
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	2H-1-Benzopyran, 3,4,4a,5,6,8a-h...	194 C13H22O	041678-32-4	38
2	Tricyclo[4.3.1.1(3,8)]undecane, ...	180 C12H20O	021898-92-0	25
3	Cyclopentanone, 2-(2-nitro-2-hep...	225 C12H19NO3	104313-57-7	25
4	3,5-Dimethyl-1-adamantanol	180 C12H20O	000707-37-9	22
5	Niacinamide, N-trimethylsilyl-	194 C9H14N2OSi	1000374-74-5	22



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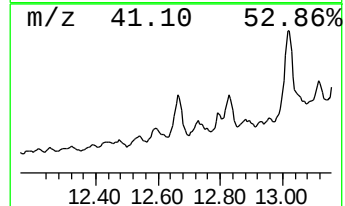
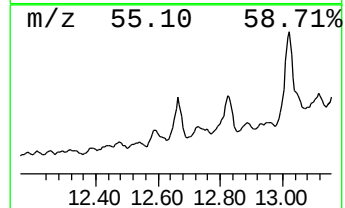
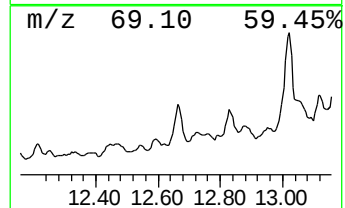
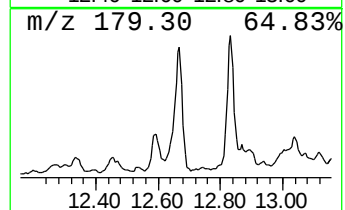
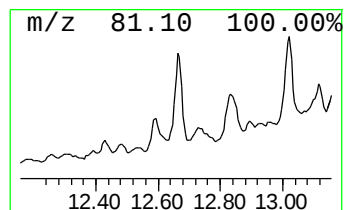
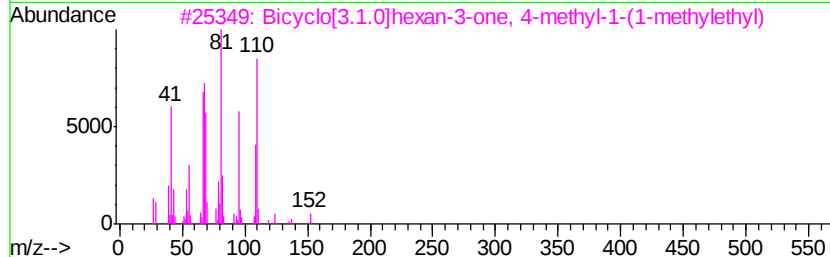
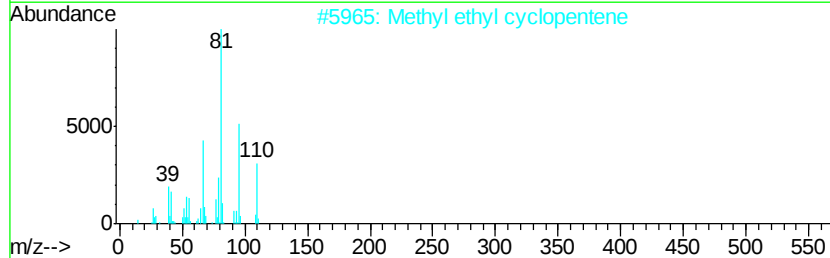
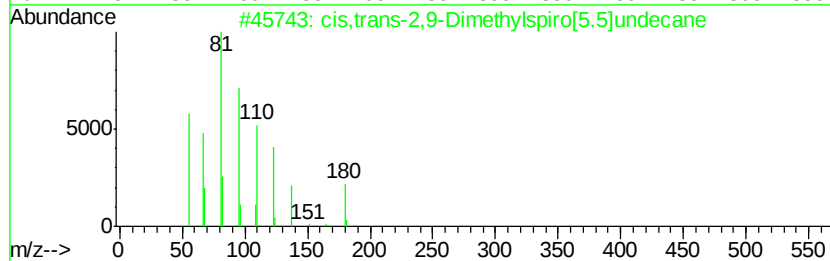
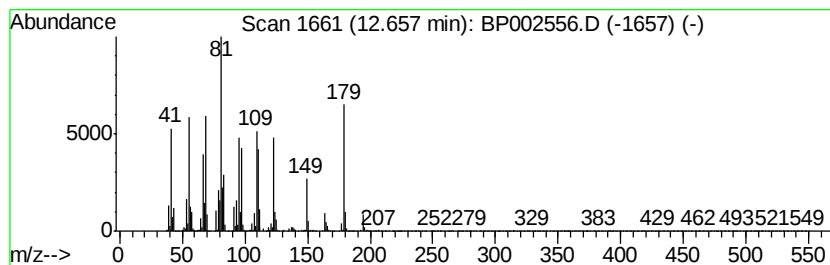
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Peak Number 3 cis,trans-2,9-Dimethylspiro... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.66	2.97 ng	994856	Acenaphthene-d10	14.24

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	cis,trans-2,9-Dimethylspiro[5.5]undecane	180	C13H24	095530-74-8	55
2		Methyl ethyl cyclopentene	110	C8H14	019780-56-4	41
3		Bicyclo[3.1.0]hexan-3-one, 4-methyl-	152	C10H16O	001125-12-8	38
4		2-Propenyl-3-vinylloxirane	110	C7H10O	070597-49-8	35
5		3-Heptyne, 5-methyl-	110	C8H14	061228-09-9	35



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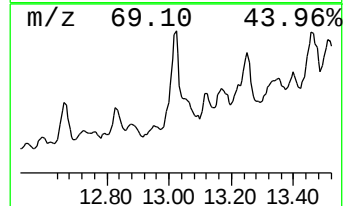
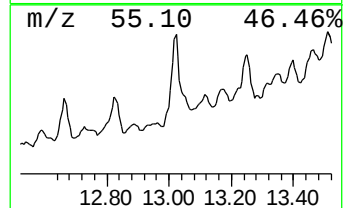
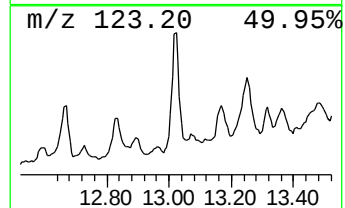
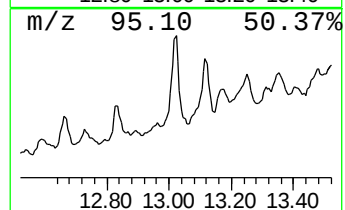
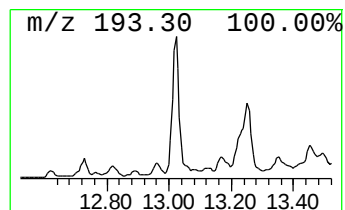
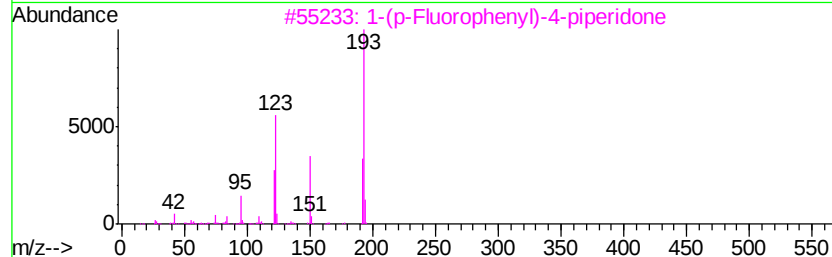
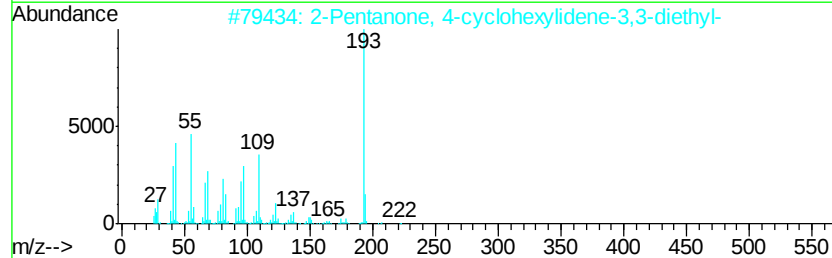
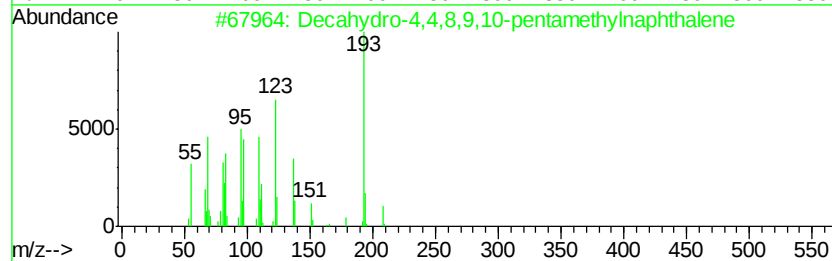
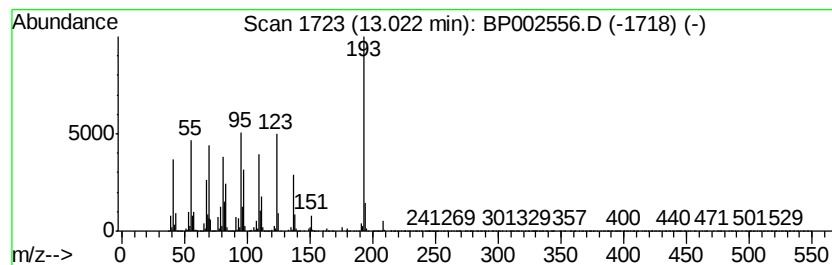
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BNA_P
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Decahydro-4,4,8,9,10-pentam... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.02	5.80 ng	1943200	Acenaphthene-d10	14.24	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Decahydro-4,4,8,9,10-pentamethyl...	208	C15H28	080655-44-3	93
2	2-Pentanone, 4-cyclohexylidene-3...	222	C15H26O	313253-65-5	50
3	1-(p-Fluorophenyl)-4-piperidone	193	C11H12FN0	1000238-56-7	47
4	2-Anthracenamine	193	C14H11N	000613-13-8	27
5	1-Anthracenamine	193	C14H11N	000610-49-1	27



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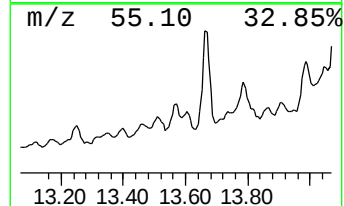
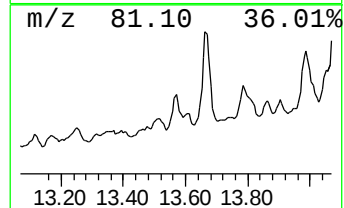
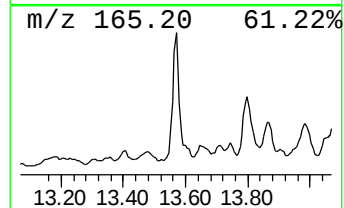
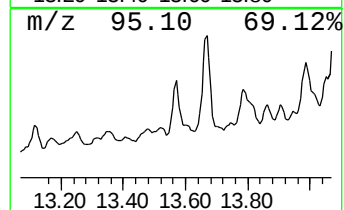
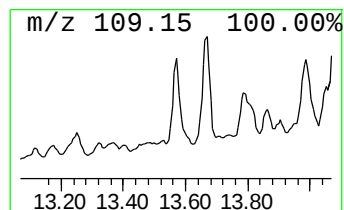
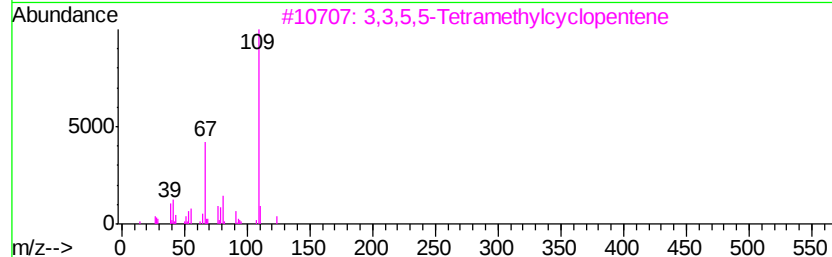
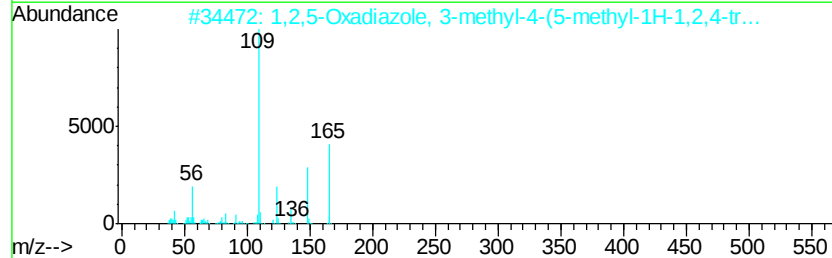
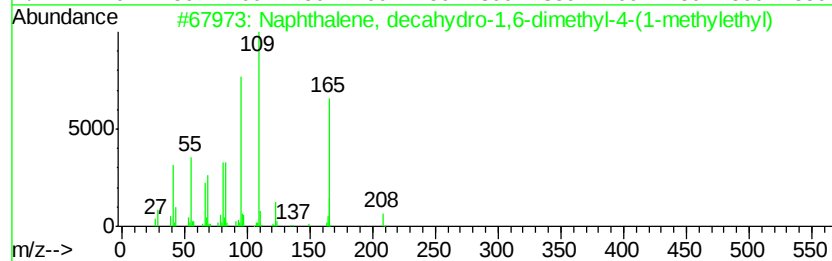
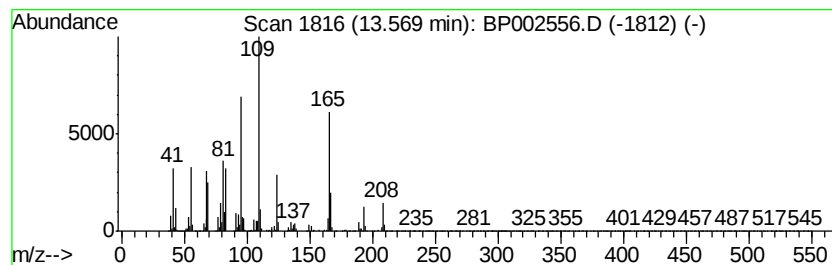
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Peak Number 5 Naphthalene, decahydro-1,6-... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.57	5.05 ng	1692580	Acenaphthene-d10	14.24

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, decahydro-1,6-dimet...	208	C15H28	029788-41-8	87
2		1,2,5-Oxadiazole, 3-methyl-4-(5-...	165	C6H7N5O	1000337-72-1	43
3		3,3,5,5-Tetramethylcyclopentene	124	C9H16	038667-10-6	30
4		4H-Chromene, 4a,5,6,7,8,8a-hexah...	208	C14H24O	1000196-77-4	30
5		2-Cyclopentene-1-carboxylic acid...	154	C9H14O2	006894-69-5	27



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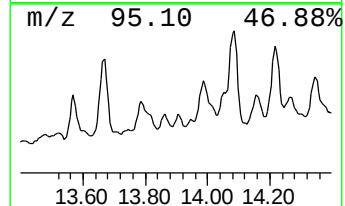
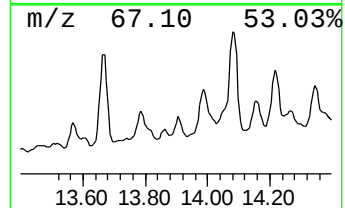
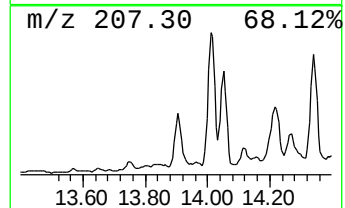
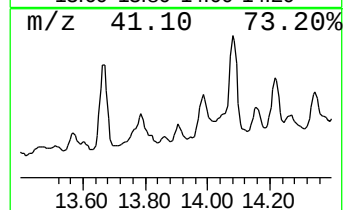
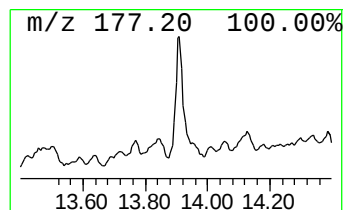
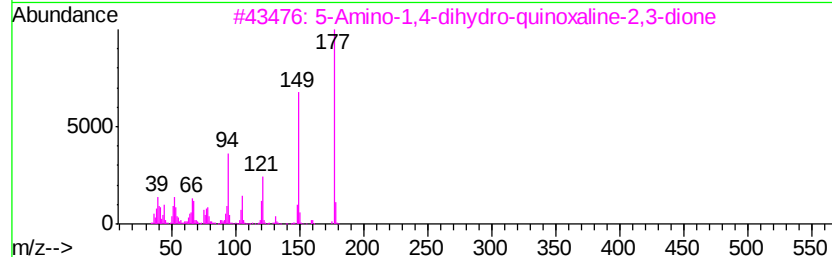
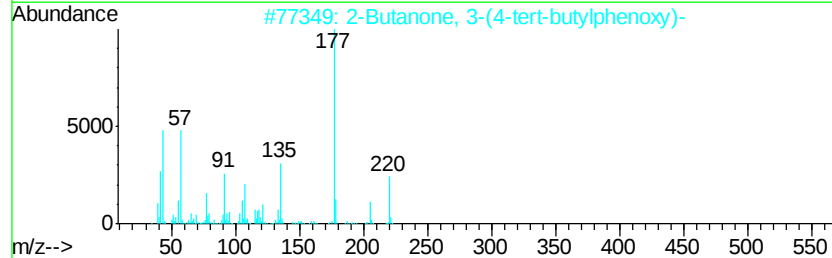
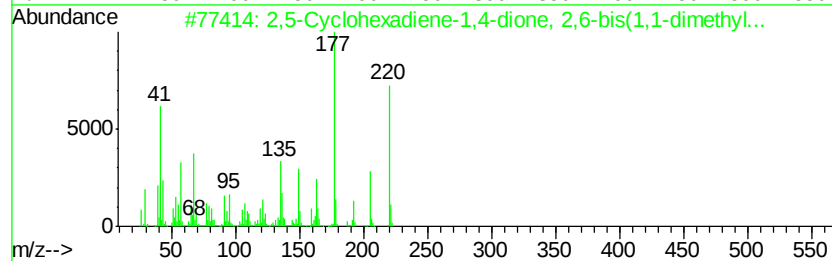
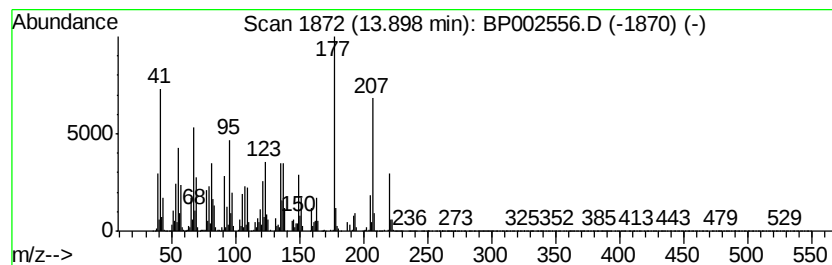
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Peak Number 6 2,5-Cyclohexadiene-1,4-dione... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.90	3.78 ng	1266240	Acenaphthene-d10	14.24

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2,5-Cyclohexadiene-1,4-dione, 2,...	220	C14H20O2	000719-22-2	90
2		2-Butanone, 3-(4-tert-butylpheno...	220	C14H20O2	160875-28-5	38
3		5-Amino-1,4-dihydro-quinoxaline-...	177	C8H7N3O2	076097-87-5	22
4		3-Buten-2-one, 4-(2,6,6-trimethy...	192	C13H20O	014901-07-6	22
5		trans-.beta.-Ionone	192	C13H20O	000079-77-6	22



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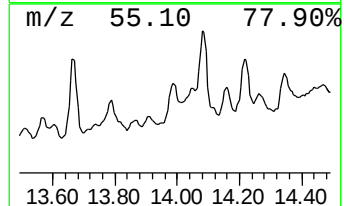
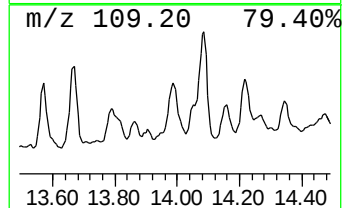
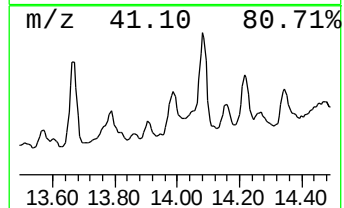
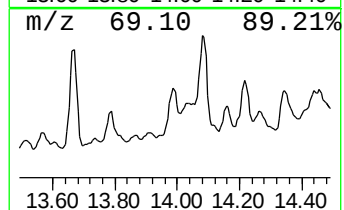
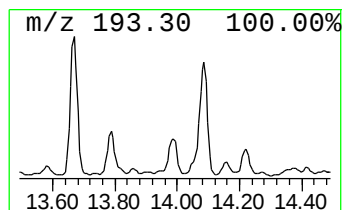
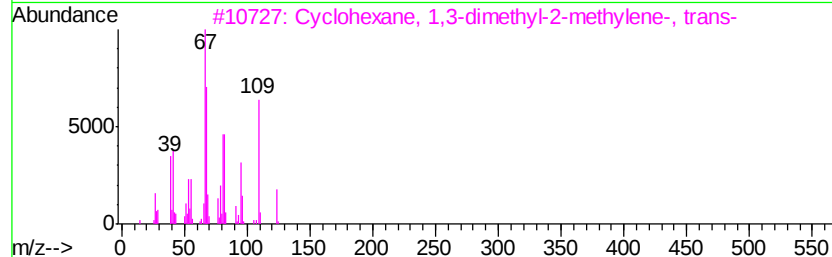
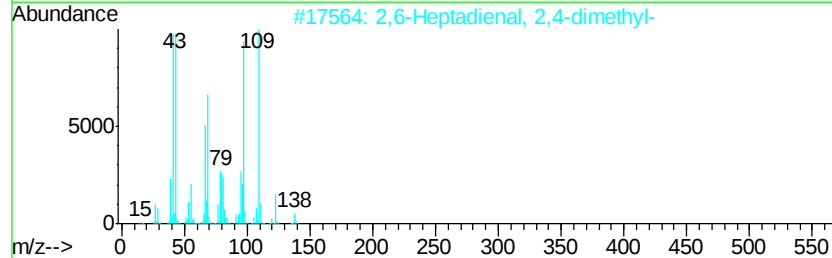
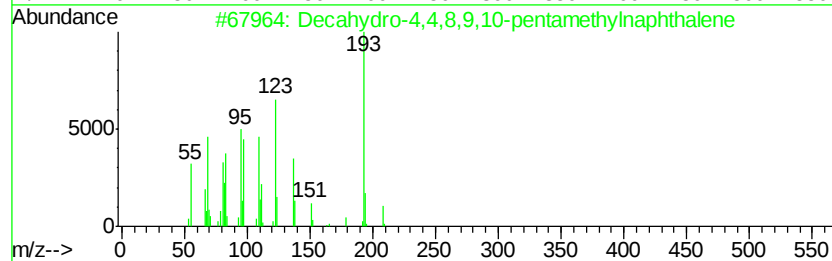
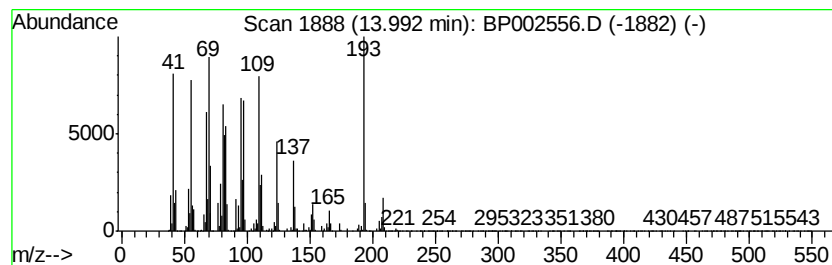
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 unknown13.99 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.99	15.01 ng	5031940	Acenaphthene-d10	14.24

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Decahydro-4,4,8,9,10-pentamethyl...	208	C15H28	080655-44-3	55
2		2,6-Heptadienal, 2,4-dimethyl-	138	C9H14O	085136-08-9	45
3		Cyclohexane, 1,3-dimethyl-2-meth...	124	C9H16	020348-74-7	35
4		Cyclopropane, 1,1-dimethyl-2-(2-...	124	C9H16	069147-03-1	27
5		3-Cyclohexene-1-carboxaldehyde, ...	152	C10H16O	040702-26-9	25



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP062620\
Data File : BP002556.D
Acq On : 26 Jun 2020 17:56
Operator : CG/JU
Sample : L3114-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

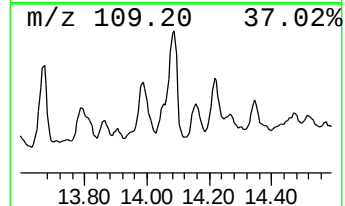
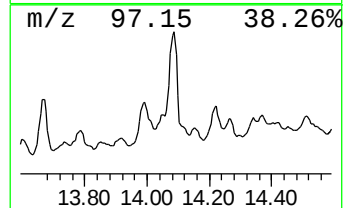
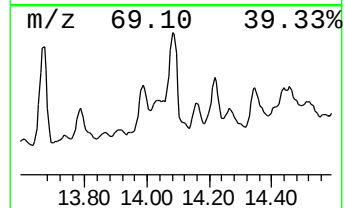
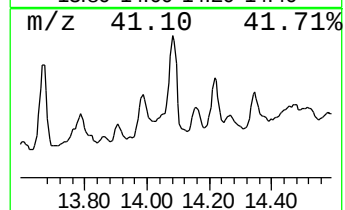
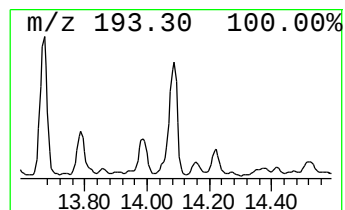
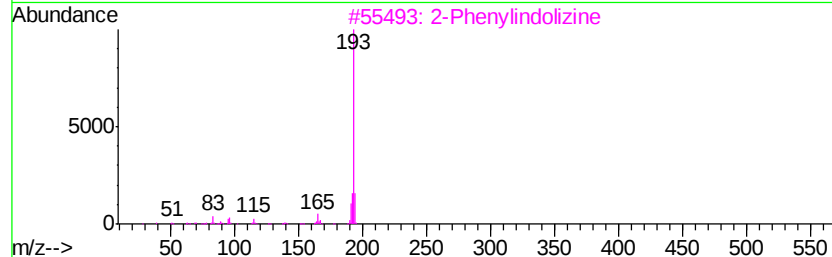
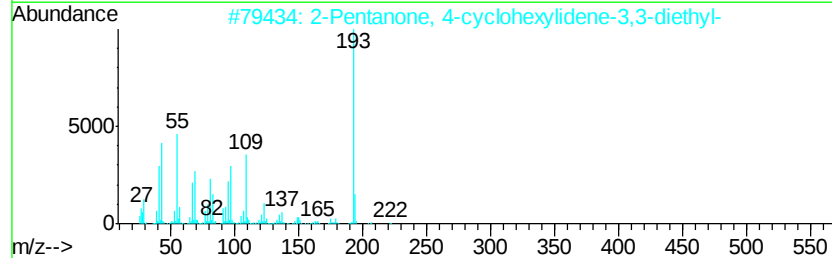
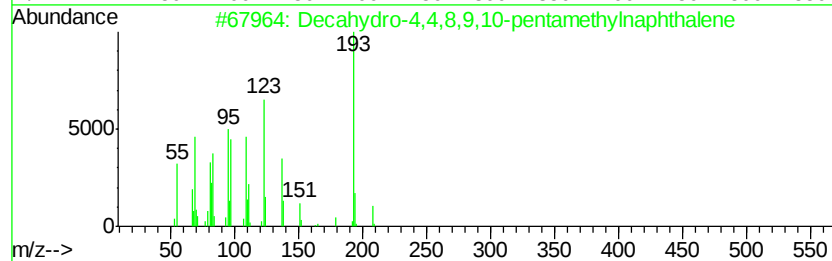
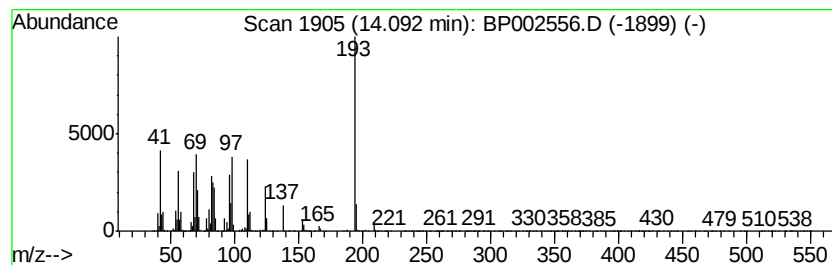
Instrument :
BNA_P
ClientSampleId :
VNJ-242

Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP060520.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 unknown14.09 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD		R.T.
14.09	22.37 ng	7500270	Acenaphthene-d10		14.24
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Decahydro-4,4,8,9,10-pentamethyl...	208	C15H28	080655-44-3	50
2	2-Pentanone, 4-cyclohexylidene-3...	222	C15H26O	313253-65-5	42
3	2-Phenylindolizine	193	C14H11N	025379-20-8	27
4	[1,1'-Biphenyl]-4-acetonitrile	193	C14H11N	031603-77-7	25
5	1H-Indene, octahydro-2,2,4,4,7,7...	208	C15H28	054832-83-6	25



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP062620\
Data File : BP002556.D
Acq On : 26 Jun 2020 17:56
Operator : CG/JU
Sample : L3114-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

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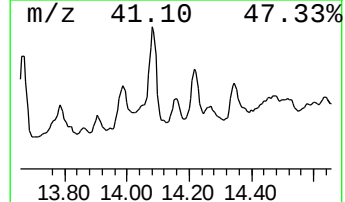
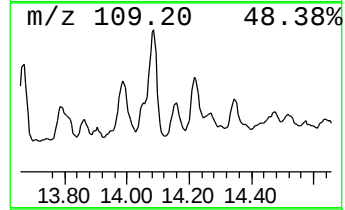
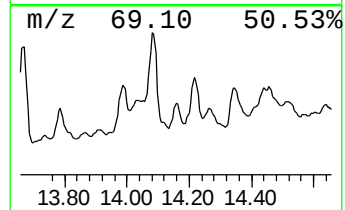
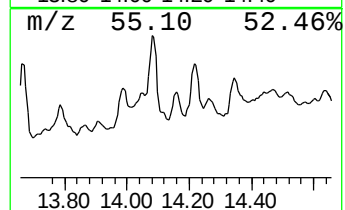
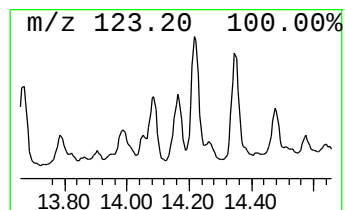
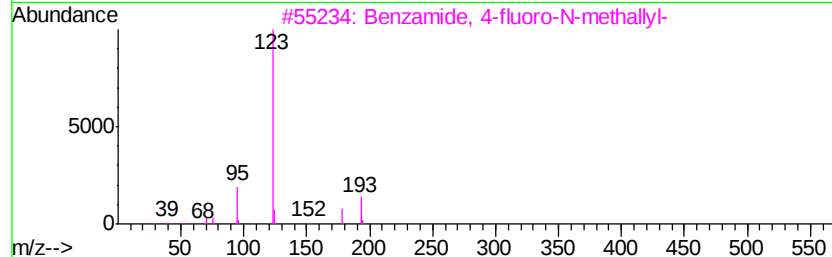
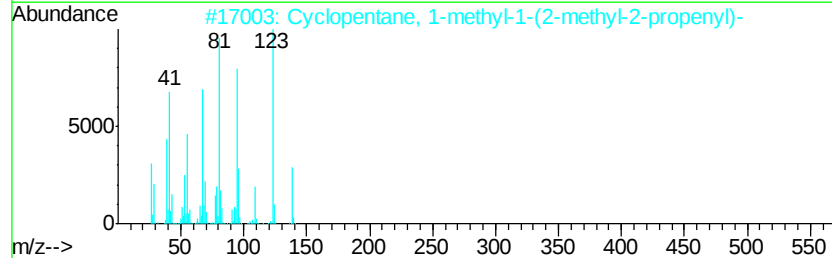
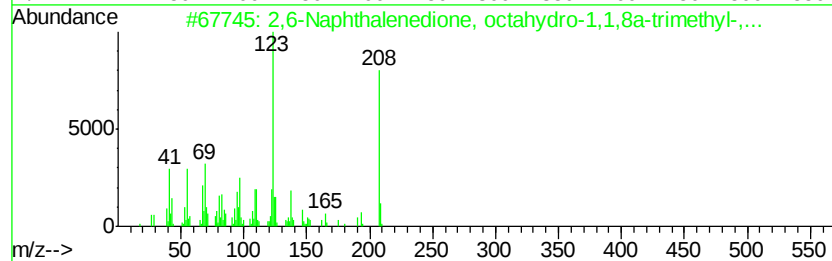
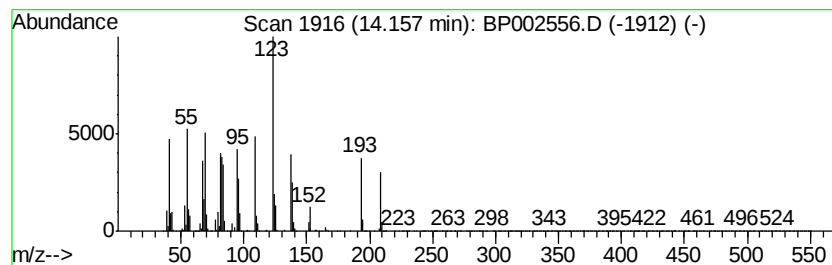
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 unknown14.16 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.16	5.57 ng	1866180	Acenaphthene-d10	14.24

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2,6-Naphthalenedione, octahydro-...	208	C13H20O2	057289-16-4	43
2			Cyclopentane, 1-methyl-1-(2-meth...	138	C10H18	074764-47-9	43
3			Benzamide, 4-fluoro-N-methyl-...	193	C11H12FNO	1000340-31-9	38
4			Cyclopentene, 1,3-dimethyl-2-(1-...	138	C10H18	061142-32-3	37
5			(2,4,6-Trimethylcyclohexyl) meth...	156	C10H20O	013702-56-2	35



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP062620\
Data File : BP002556.D
Acq On : 26 Jun 2020 17:56
Operator : CG/JU
Sample : L3114-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

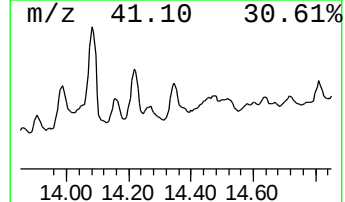
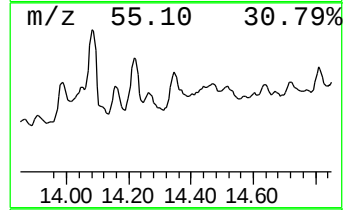
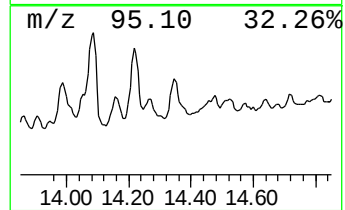
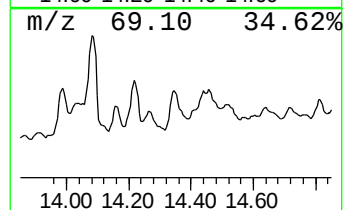
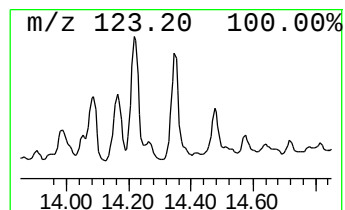
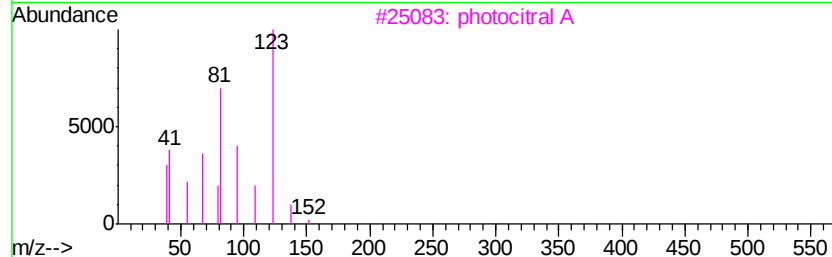
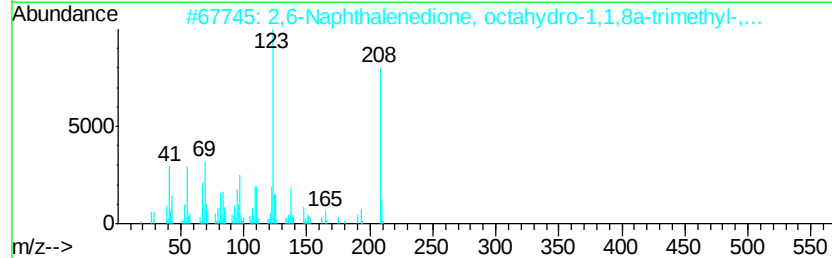
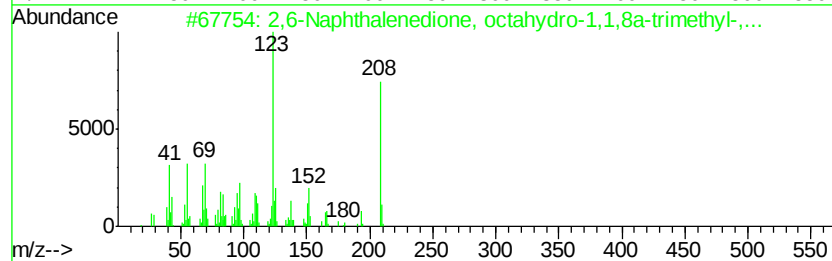
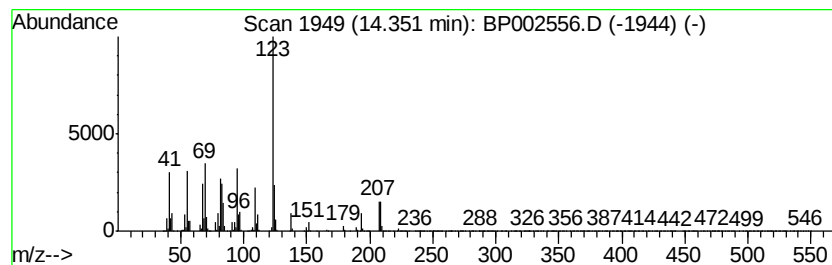
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP060520.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 2,6-Naphthalenedione, octah... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.35	10.03 ng	3362020	Acenaphthene-d10	14.24
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	2,6-Naphthalenedione, octahydro-...	208 C13H20O2	057289-17-5	64
2	2,6-Naphthalenedione, octahydro-...	208 C13H20O2	057289-16-4	50
3	photocitral A	152 C10H16O	055253-28-6	43
4	Cyclohexene, 1,5,5-trimethyl-6-a...	180 C12H20O	211563-96-1	42
5	4-Fluorobenzoic acid, 4-pentadec...	350 C22H35FO2	1000280-68-5	38



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP062620\
Data File : BP002556.D
Acq On : 26 Jun 2020 17:56
Operator : CG/JU
Sample : L3114-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

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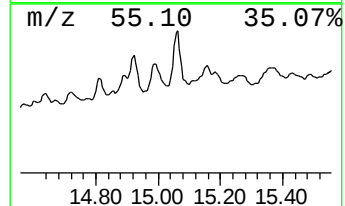
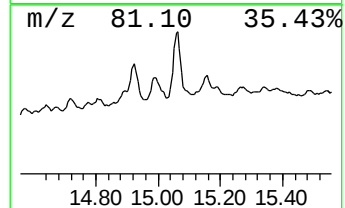
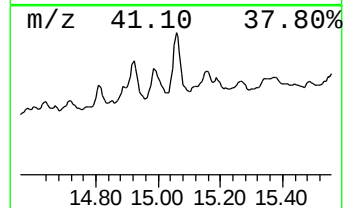
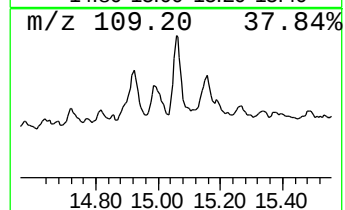
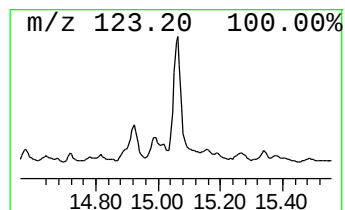
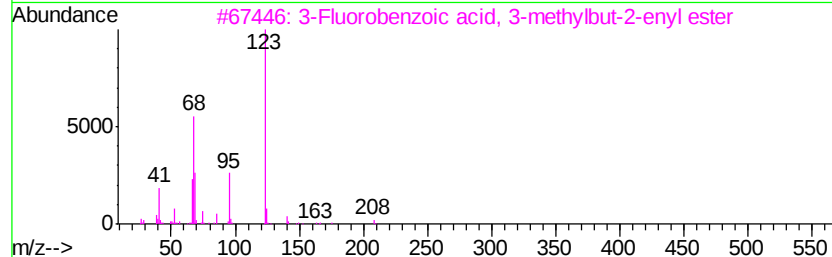
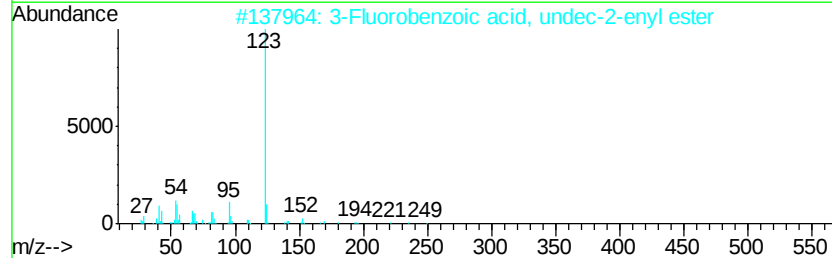
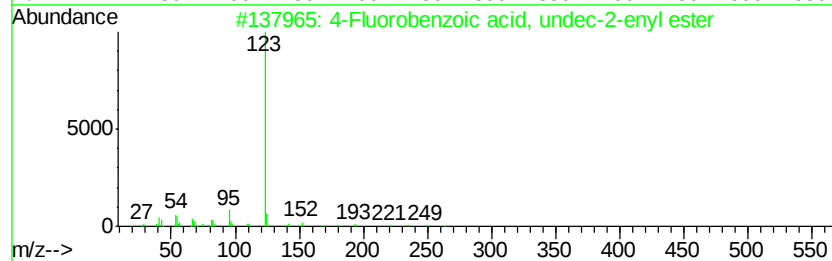
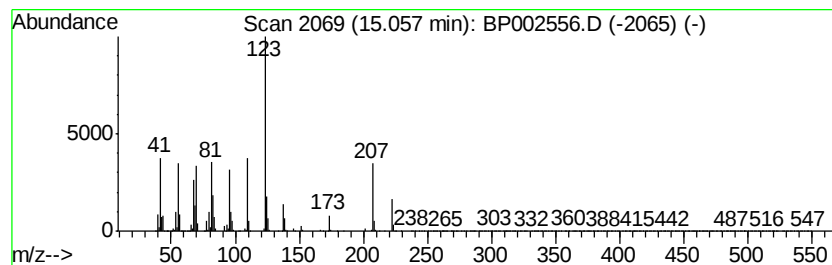
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 4-Fluorobenzoic acid, undec... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.06	13.27 ng	4449010	Acenaphthene-d10	14.24

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4-Fluorobenzoic acid, undec-2-en...	292	C18H25F02	1000299-15-8	50
2		3-Fluorobenzoic acid, undec-2-en...	292	C18H25F02	1000292-60-8	50
3		3-Fluorobenzoic acid, 3-methylbu...	208	C12H13F02	154559-38-3	47
4		2-Fluorobenzoic acid, 2-ethylcyc...	250	C15H19F02	1000278-98-1	43
5		Cyclohexene, 1,5,5-trimethyl-6-a...	180	C12H200	211563-96-1	40



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP062620\
Data File : BP002556.D
Acq On : 26 Jun 2020 17:56
Operator : CG/JU
Sample : L3114-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

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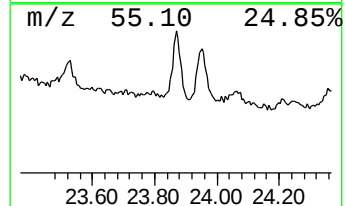
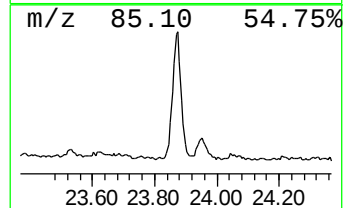
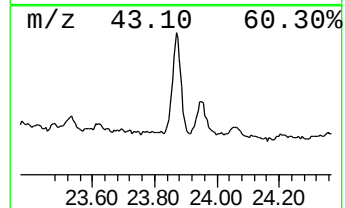
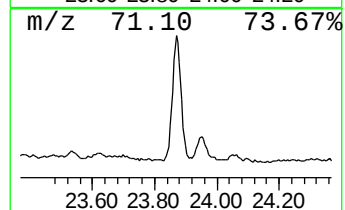
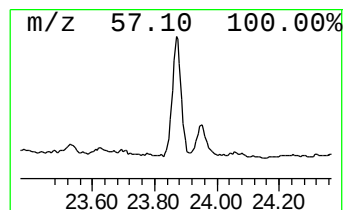
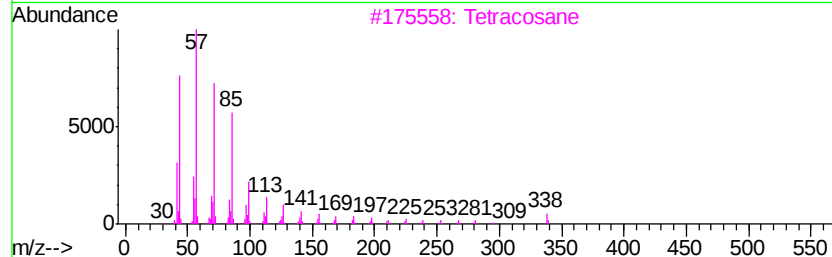
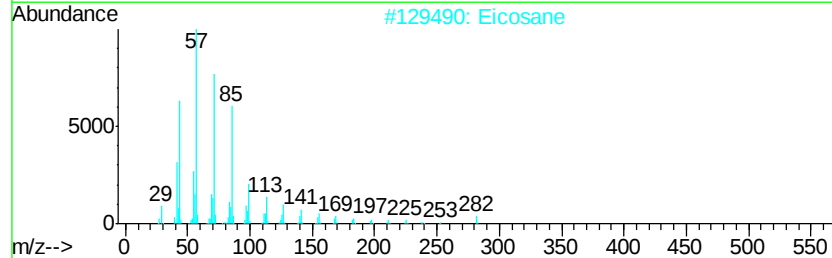
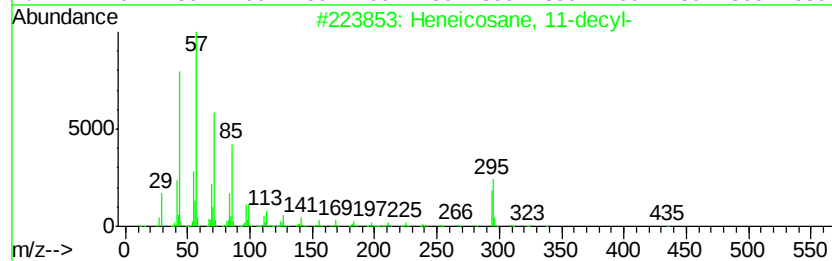
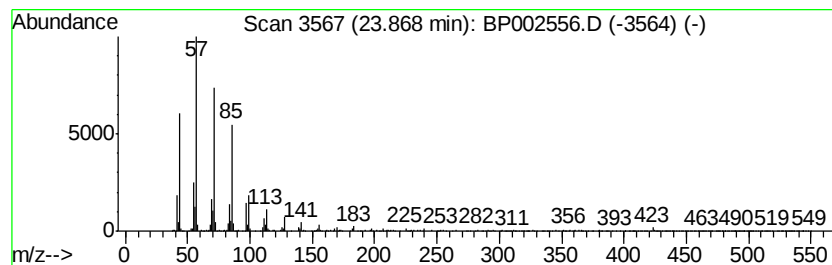
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Heneicosane, 11-decyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.87	8.36 ng	579274	Perylene-d12	23.30

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heneicosane, 11-decyl-	437	C31H64	055320-06-4	95
2			Eicosane	282	C20H42	000112-95-8	94
3			Tetracosane	338	C24H50	000646-31-1	94
4			Octacosane	394	C28H58	000630-02-4	91
5			Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP062620\
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 Acq On : 26 Jun 2020 17:56
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 Sample : L3114-01
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Instrument :
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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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
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cis,trans-2,9-Dim...	12.66	3.0	ng	994856	3	14.24	6704830	20.0
Decahydro-4,4,8,9...	13.02	5.8	ng	1943200	3	14.24	6704830	20.0
Naphthalene, deca...	13.57	5.0	ng	1692580	3	14.24	6704830	20.0
2,5-Cyclohexadien...	13.90	3.8	ng	1266240	3	14.24	6704830	20.0
unknown13.99	13.99	15.0	ng	5031940	3	14.24	6704830	20.0
unknown14.09	14.09	22.4	ng	7500270	3	14.24	6704830	20.0
unknown14.16	14.16	5.6	ng	1866180	3	14.24	6704830	20.0
2,6-Naphthalenedi...	14.35	10.0	ng	3362020	3	14.24	6704830	20.0
4-Fluorobenzoic a...	15.06	13.3	ng	4449010	3	14.24	6704830	20.0
Heneicosane, 11-d...	23.87	8.4	ng	579274	6	23.30	1385000	20.0