

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051321\
 Data File : BP005530.D
 Acq On : 13 May 2021 21:09
 Operator : CG/JU
 Sample : M2372-09
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 CHRT-26516

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_P\Methods\8270E-BP050521.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BP005530.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.281	366	373	389	rBV	269245	391807	26.79%	5.673%
2	6.887	638	646	662	rBV	229898	385472	26.36%	5.581%
3	7.687	775	782	790	rBV	80707	124049	8.48%	1.796%
4	8.857	971	981	988	rBV	158639	273079	18.67%	3.954%
5	10.475	1249	1256	1264	rBV	88301	152470	10.43%	2.208%
6	12.757	1640	1644	1650	rVB6	13302	19364	1.32%	0.280%
7	12.957	1667	1678	1686	rBV	547586	775389	53.02%	11.227%
8	13.098	1698	1702	1710	rBV2	31264	50633	3.46%	0.733%
9	13.745	1807	1812	1816	rBV	52376	77829	5.32%	1.127%
10	13.810	1818	1823	1828	rVV2	52818	94015	6.43%	1.361%
11	13.863	1828	1832	1841	rVB8	28408	62349	4.26%	0.903%
12	13.975	1848	1851	1856	rBV6	15244	25906	1.77%	0.375%
13	14.063	1862	1866	1871	rBV4	41187	64472	4.41%	0.933%
14	14.116	1872	1875	1877	rVV3	28522	33536	2.29%	0.486%
15	14.151	1877	1881	1888	rVB2	90820	142706	9.76%	2.066%
16	14.234	1890	1895	1898	rBV6	29142	54792	3.75%	0.793%
17	14.287	1900	1904	1908	rVV3	48731	84143	5.75%	1.218%
18	14.340	1908	1913	1920	rVB	176929	268507	18.36%	3.888%
19	14.863	1999	2002	2007	rVB4	47682	62158	4.25%	0.900%
20	14.987	2015	2023	2029	rBV9	50561	116371	7.96%	1.685%
21	15.051	2029	2034	2040	rBV9	48535	120580	8.25%	1.746%
22	15.116	2041	2045	2050	rBV2	144392	209117	14.30%	3.028%
23	15.839	2164	2168	2174	rBV	621801	810287	55.41%	11.732%
24	17.098	2379	2382	2388	rBV	309223	494398	33.81%	7.158%
25	19.669	2815	2819	2828	rVB	1168973	1462372	100.00%	21.174%
26	21.192	3074	3078	3083	rBV	216863	280830	19.20%	4.066%
27	23.504	3465	3471	3477	rVB2	146746	269948	18.46%	3.909%

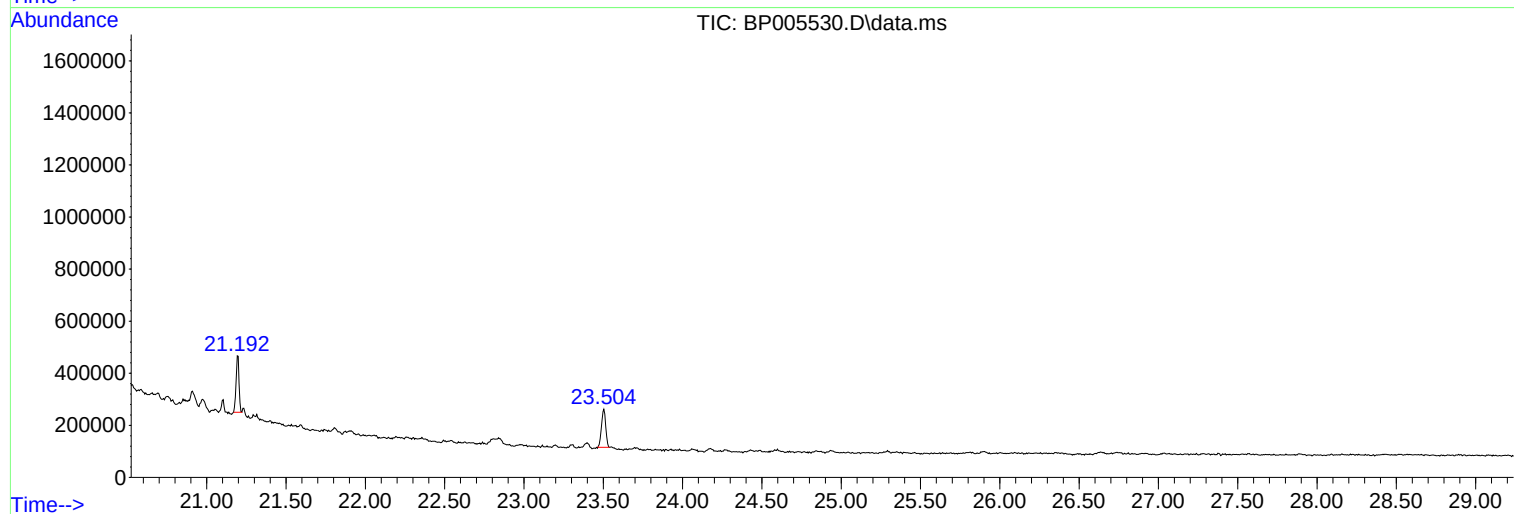
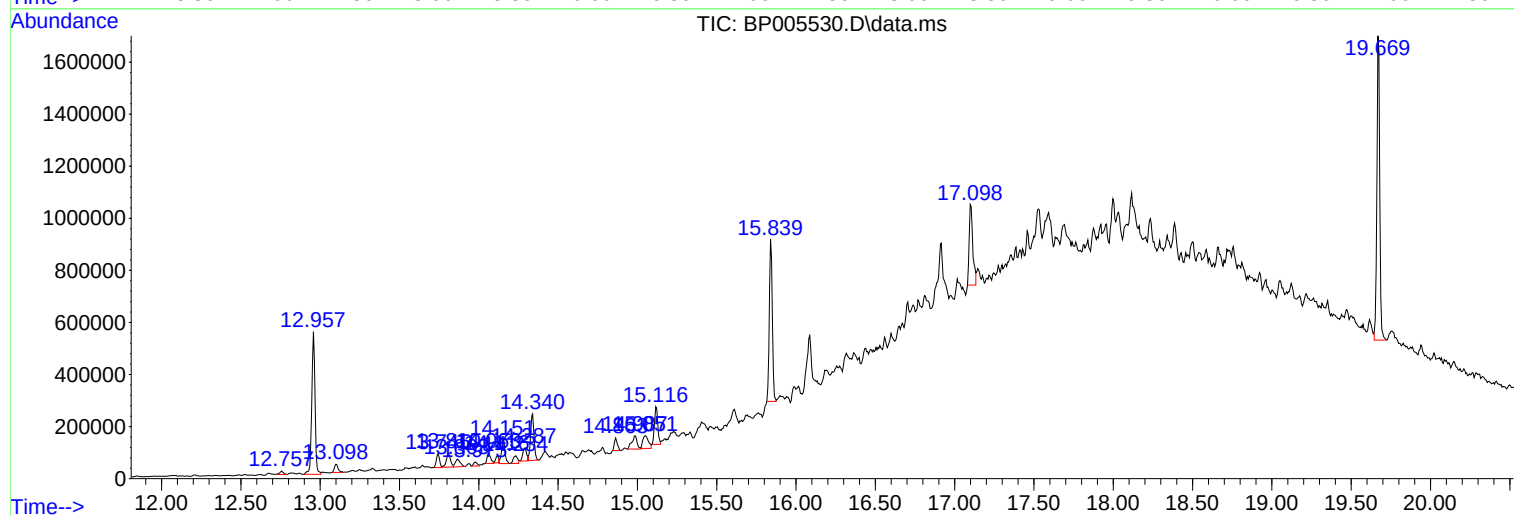
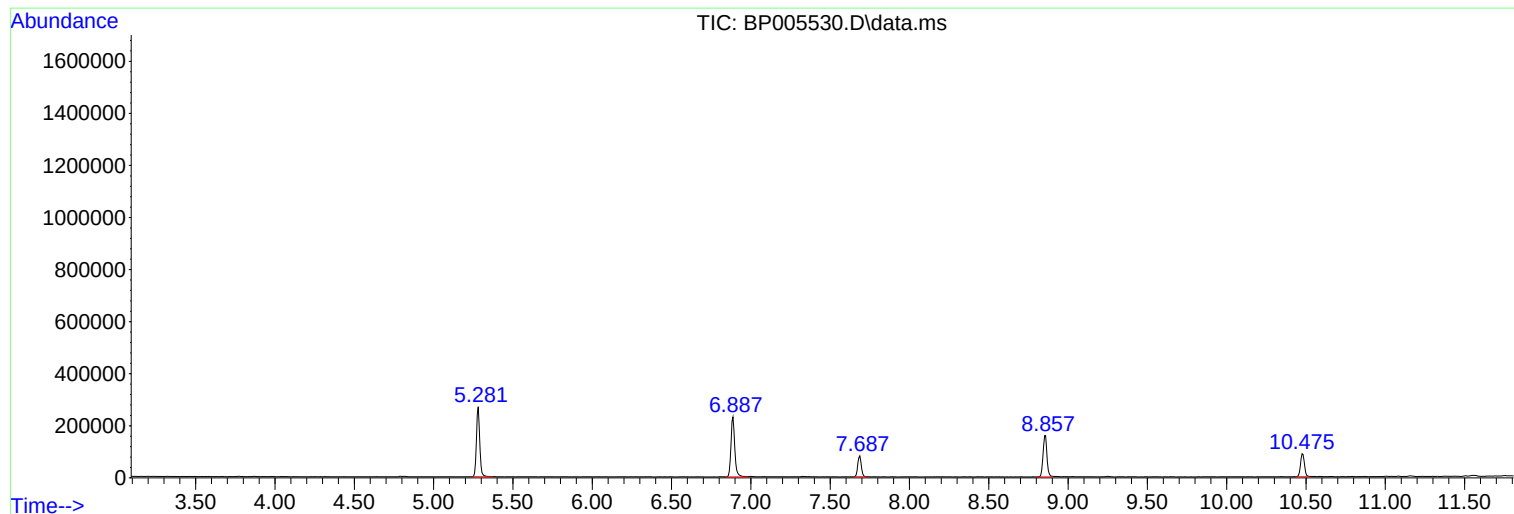
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Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP050521.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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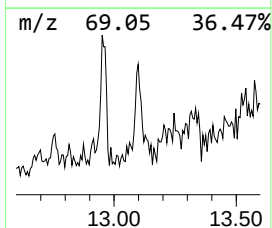
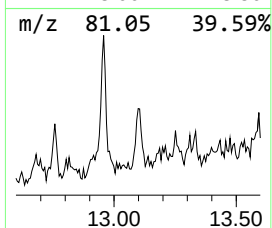
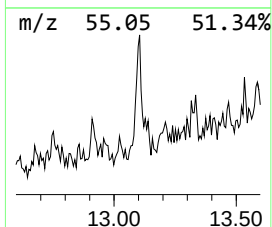
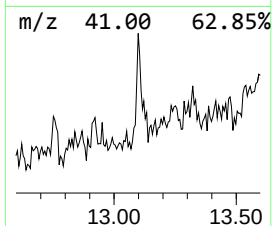
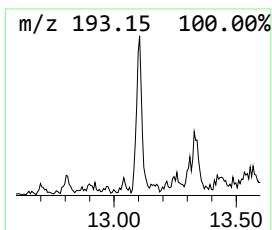
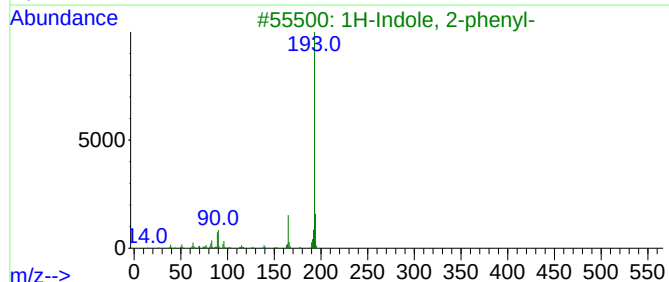
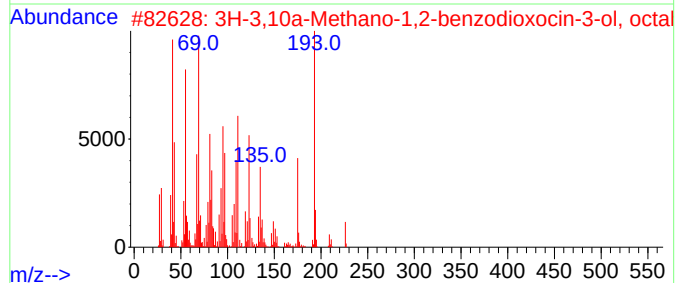
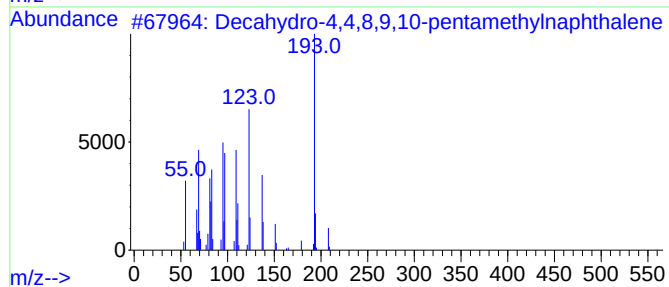
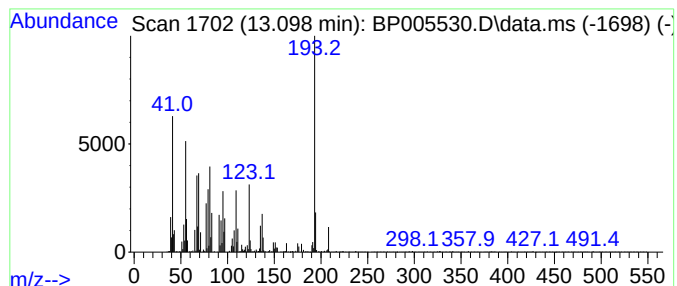
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Decahydro-4,4,8,9,10-pentam... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.098	3.77 ng	50633	Acenaphthene-d10	14.339

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Decahydro-4,4,8,9,10-pentamethyl...	208	C15H28	080655-44-3	91
2			3H-3,10a-Methano-1,2-benzodioxoc...	226	C13H22O3	095906-83-5	50
3			1H-Indole, 2-phenyl-	193	C14H11N	000948-65-2	43
4			3-Phenylindole	193	C14H11N	001504-16-1	43
5			Iminostilbene	193	C14H11N	000256-96-2	43



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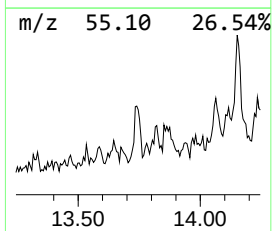
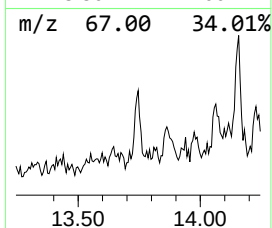
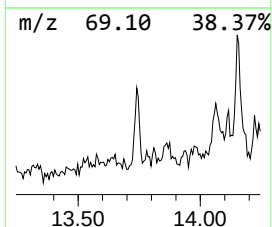
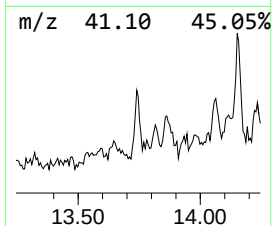
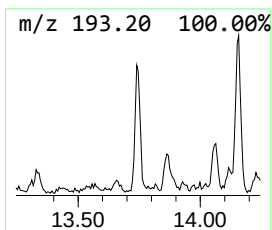
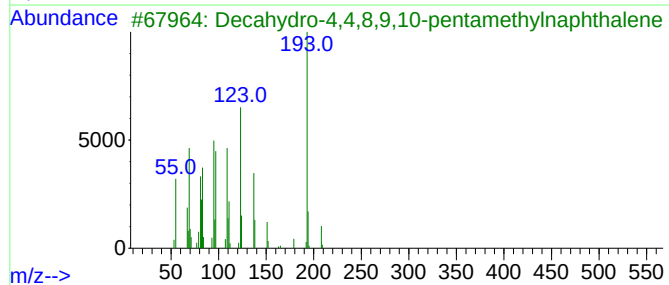
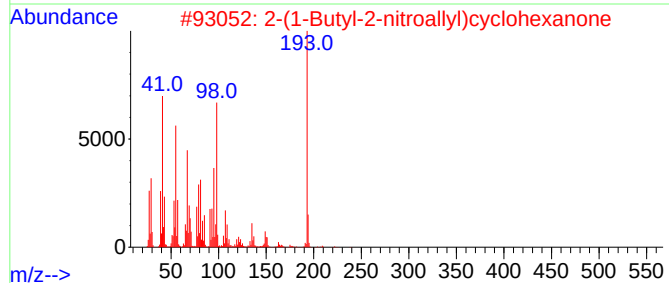
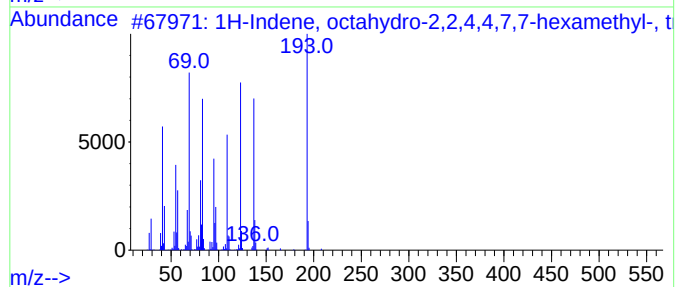
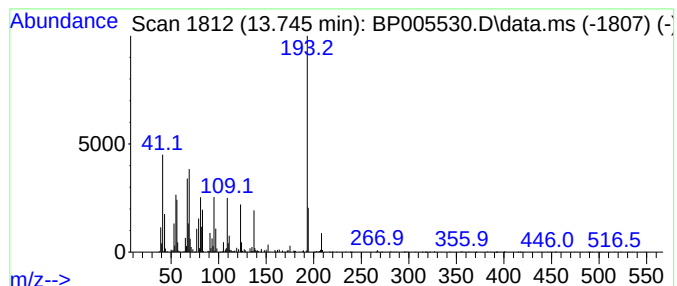
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 1H-Indene, octahydro-2,2,4,... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.746	5.80 ng	77829	Acenaphthene-d10	14.339

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Indene, octahydro-2,2,4,4,7,7...	208	C15H28	054832-83-6	59
2			2-(1-Butyl-2-nitroallyl)cyclohex...	239	C13H21NO3	1000192-05-9	38
3			Decahydro-4,4,8,9,10-pentamethyl...	208	C15H28	080655-44-3	35
4			1,5-Hexadien-3-ol, 3-methyl-6-(m...	280	C17H28OS	097369-78-3	33
5			1-(2-Hydroxyethyl)-1,2,5,5-tetra...	238	C16H30O	1000298-97-4	25



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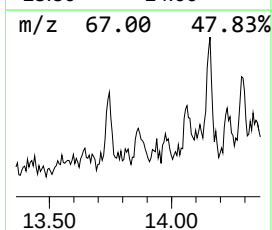
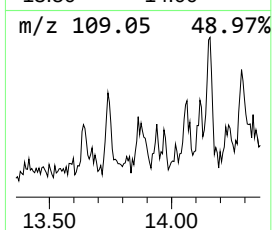
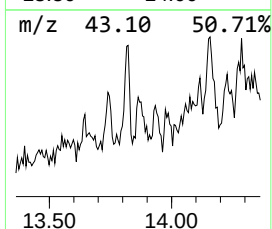
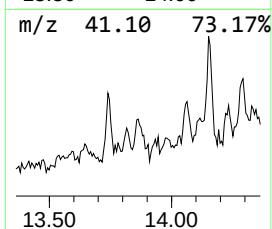
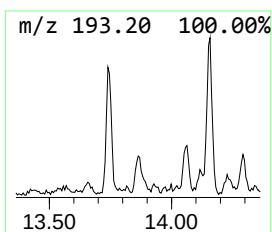
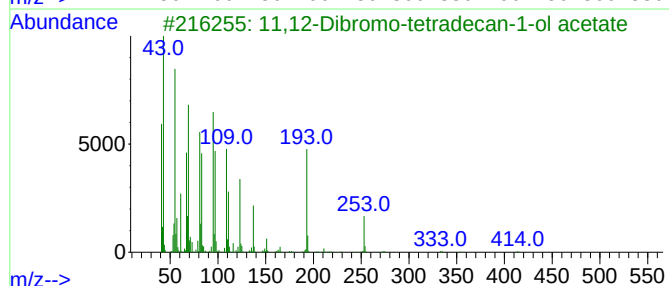
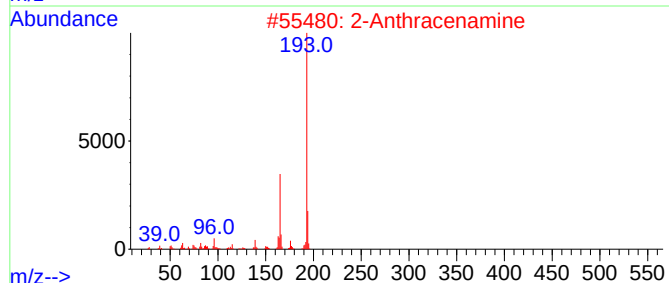
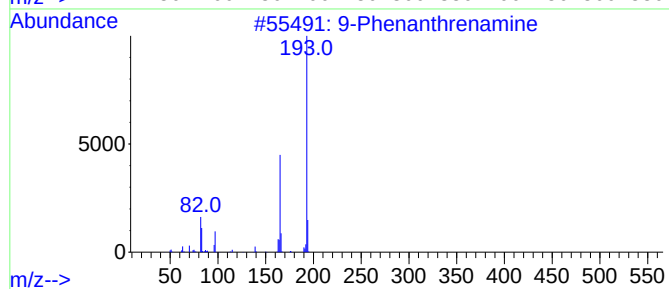
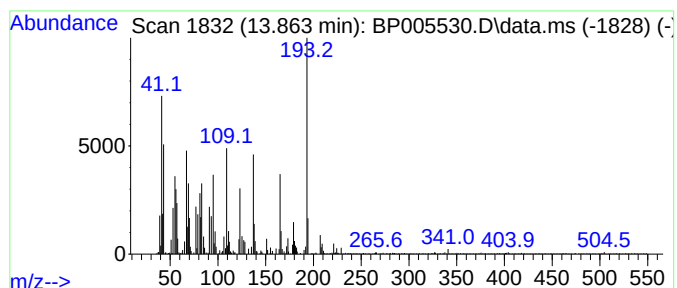
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 unknown13.863 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.863	4.64 ng	62349	Acenaphthene-d10	14.339

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			9-Phenanthrenamine	193	C14H11N	000947-73-9	42
2			2-Anthracenamine	193	C14H11N	000613-13-8	42
3			11,12-Dibromo-tetradecan-1-ol ac...	412	C16H30Br2O2	1000130-78-5	32
4			2-Biphenylacetoneitrile	193	C14H11N	019853-10-2	27
5			3-Phenylindole	193	C14H11N	001504-16-1	27



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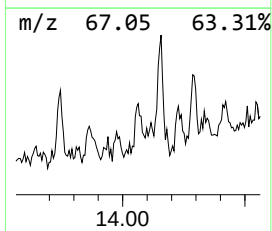
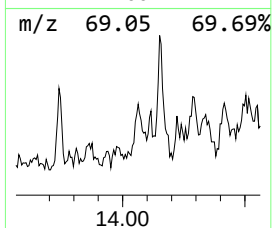
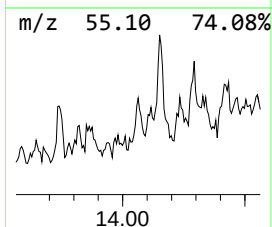
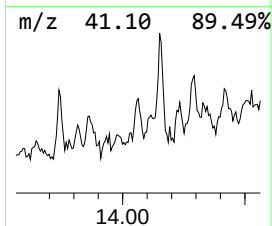
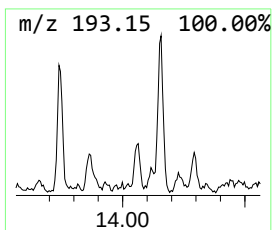
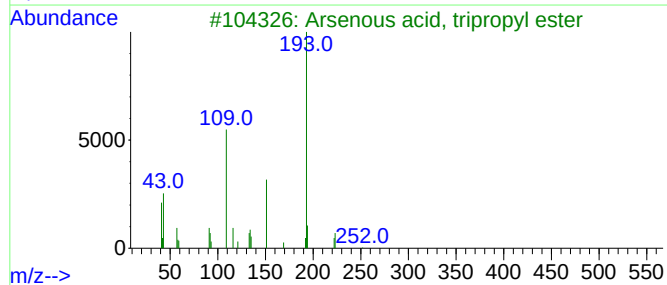
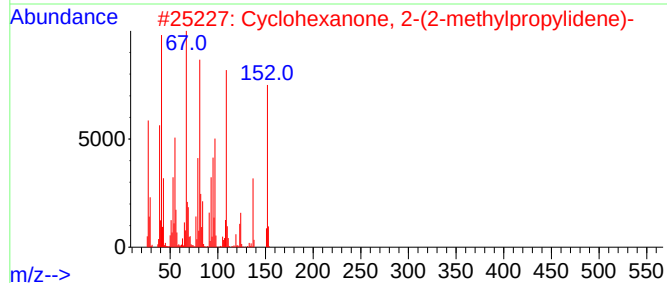
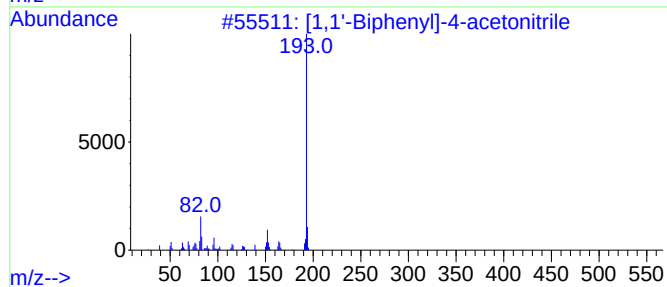
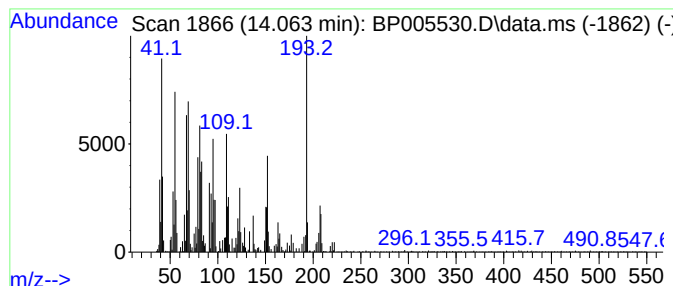
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 4 unknown14.063 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.063	4.80 ng	64472	Acenaphthene-d10	14.339

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			[1,1'-Biphenyl]-4-acetonitrile	193	C14H11N	031603-77-7	45
2			Cyclohexanone, 2-(2-methylpropylidene)-	152	C10H16O	043108-69-6	38
3			Arsenous acid, tripropyl ester	252	C9H21AsO3	015606-91-4	38
4			Naphthalene, decahydro-2-methyl-	152	C11H20	002958-76-1	35
5			1-Methyldecahydronaphthalene	152	C11H20	002958-75-0	35



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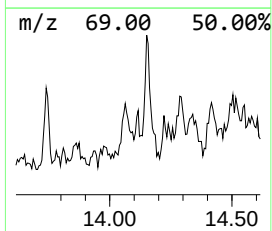
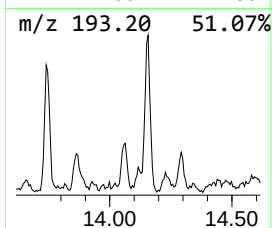
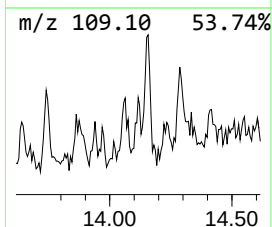
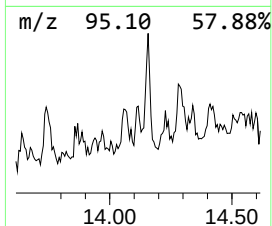
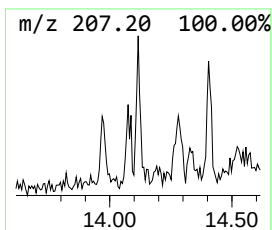
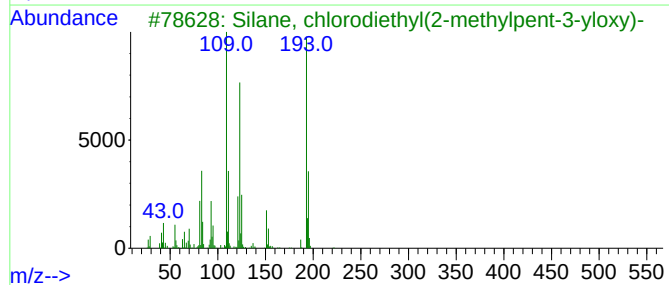
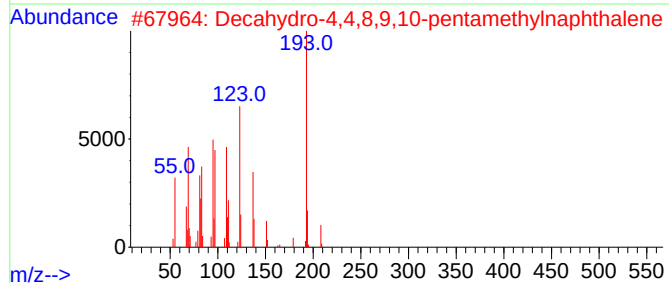
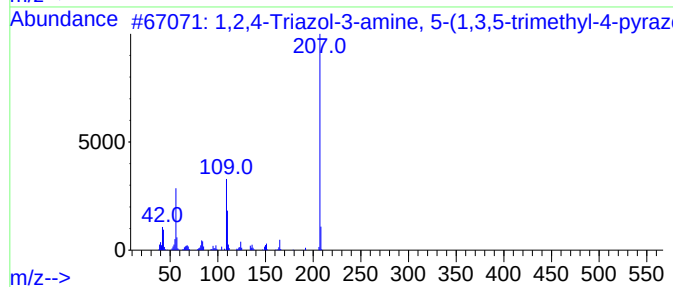
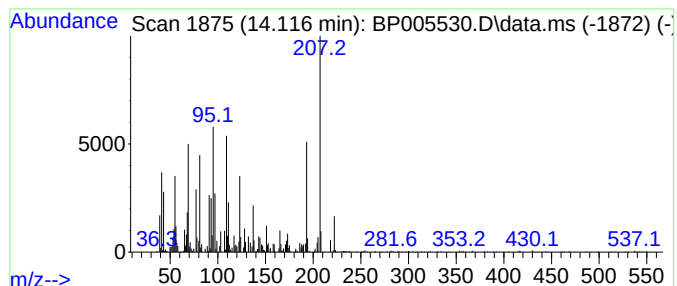
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 5 unknown14.116 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.116	2.50 ng	33536	Acenaphthene-d10	14.339

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,2,4-Triazol-3-amine, 5-(1,3,5-...	207	C8H13N7	1000264-16-7	25		
2	Decahydro-4,4,8,9,10-pentamethyl...	208	C15H28	080655-44-3	22		
3	Silane, chlorodiethyl(2-methylpe...	222	C10H23ClOSi	1000363-79-1	16		
4	Thieno[2,3-c]furan-3-carbonitril...	222	C11H14N2OS	447412-24-0	14		
5	1,2,4-Benznetricarboxylic acid, ...	224	C10H8O6	1000257-89-4	14		



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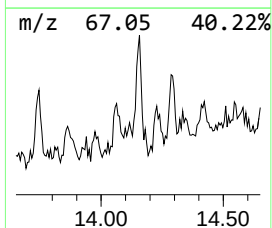
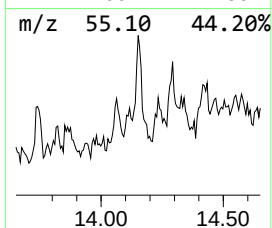
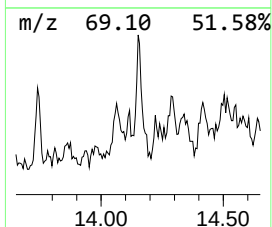
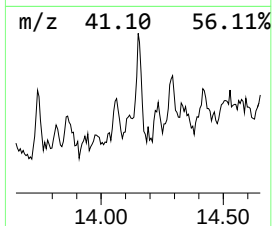
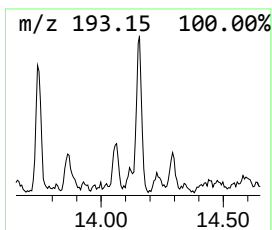
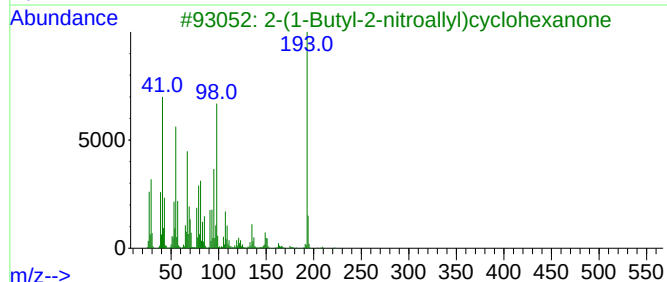
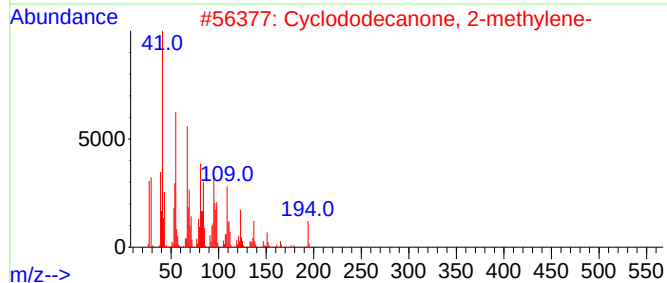
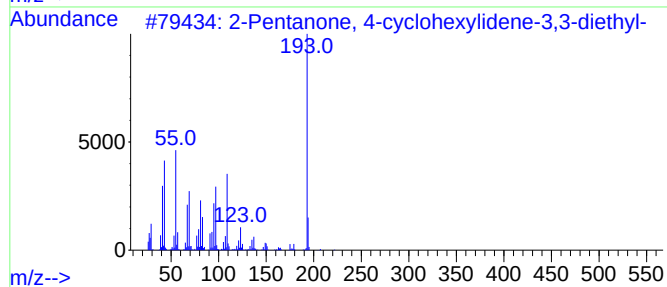
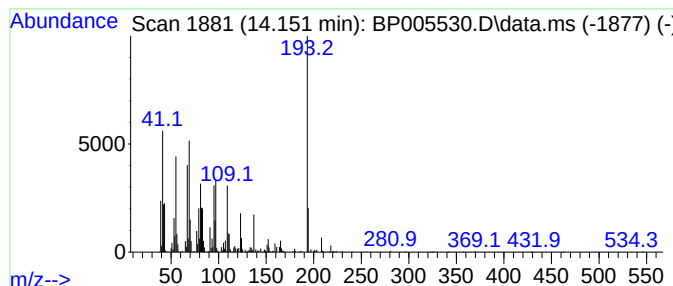
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 2-Pentanone, 4-cyclohexylid... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.151	10.63 ng	142706	Acenaphthene-d10	14.339

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-cyclohexylidene-3...	222	C15H26O	313253-65-5	58		
2	Cyclododecanone, 2-methylene-	194	C13H22O	003045-76-9	47		
3	2-(1-Butyl-2-nitroallyl)cyclohex...	239	C13H21NO3	1000192-05-9	42		
4	2-Diethylamino-4-phenylthiooct-2...	302	C18H26N2S	1000090-57-0	35		
5	6-Methylphenanthridine	193	C14H11N	003955-65-5	32		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051321\
Data File : BP005530.D
Acq On : 13 May 2021 21:09
Operator : CG/JU
Sample : M2372-09
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CHRT-26516

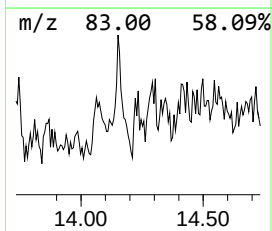
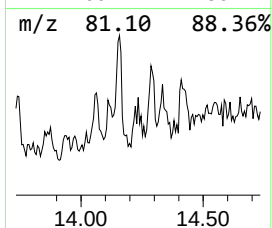
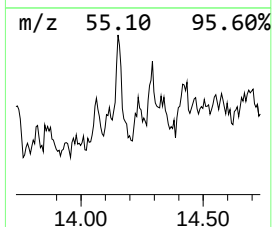
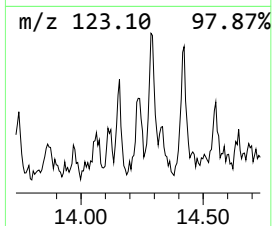
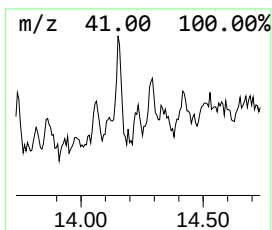
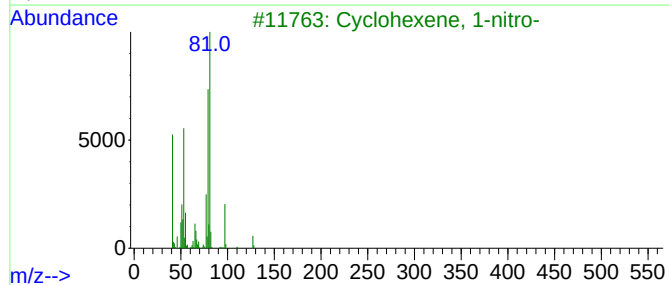
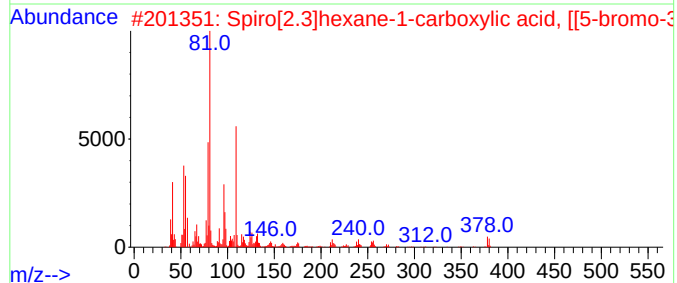
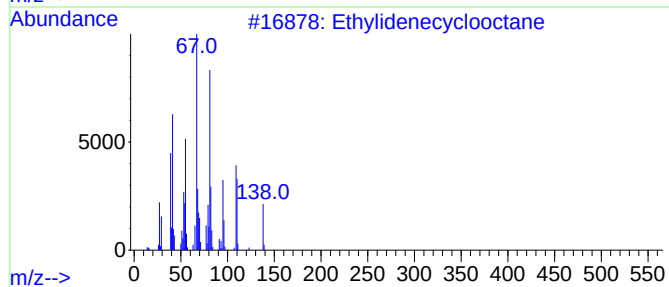
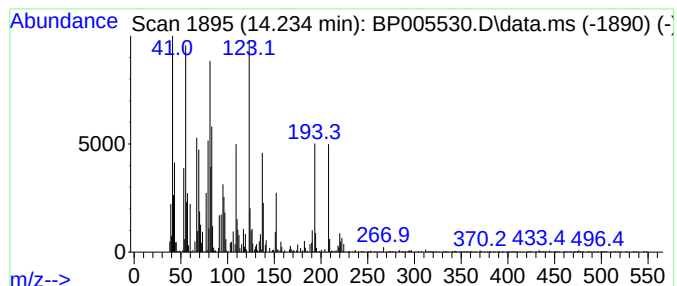
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 unknown14.234 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.234	4.08 ng	54792	Acenaphthene-d10	14.339

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethylidenecyclooctane	138	C10H18	019780-51-9	43
2			Spiro[2.3]hexane-1-carboxylic ac...	378	C18H23BrN2O2	330966-62-6	38
3			Cyclohexene, 1-nitro-	127	C6H9NO2	002562-37-0	38
4			2-Hexyne, 6-iodo-	208	C6H9I	028077-74-9	38
5			1-Cyclohexene-1-methanol	112	C7H12O	004845-04-9	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051321\
Data File : BP005530.D
Acq On : 13 May 2021 21:09
Operator : CG/JU
Sample : M2372-09
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CHRT-26516

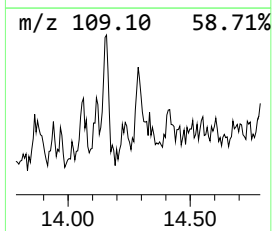
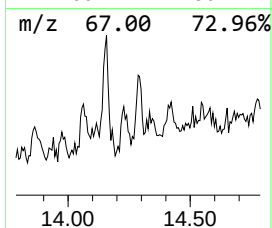
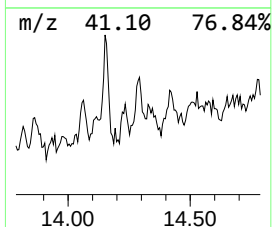
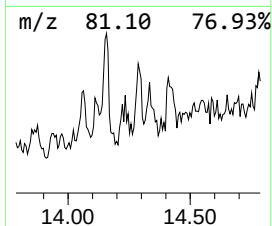
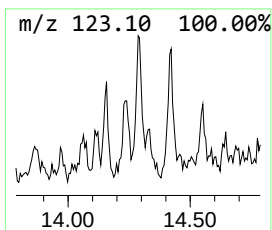
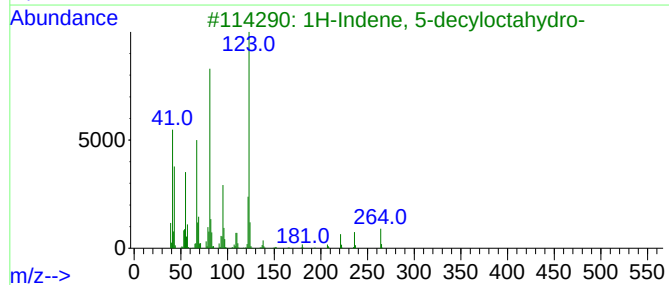
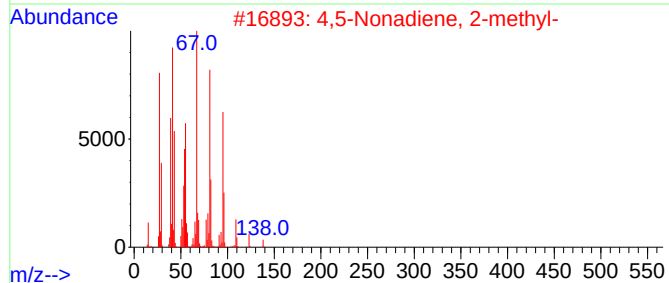
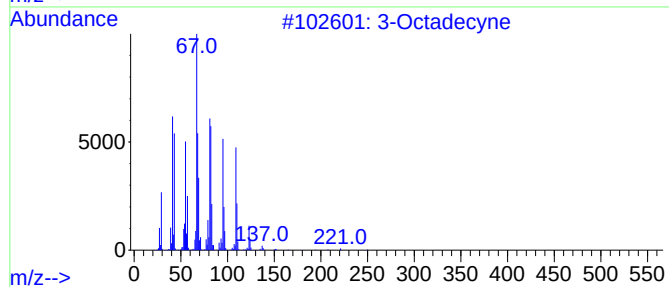
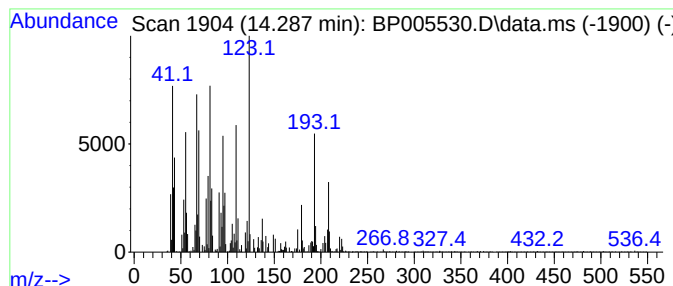
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 unknown14.287 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.287	6.27 ng	84143	Acenaphthene-d10	14.339

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Octadecyne	250	C18H34	061886-64-4	43
2			4,5-Nonadiene, 2-methyl-	138	C10H18	055956-32-6	42
3			1H-Indene, 5-decyloctahydro-	264	C19H36	055044-35-4	38
4			1-(Phenylsulfonylamino)-5-methyl-...	292	C13H16N4O2S	120675-04-9	38
5			Bicyclo[2.2.1]heptane, 1,3,3-tri...	138	C10H18	006248-88-0	27



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051321\
Data File : BP005530.D
Acq On : 13 May 2021 21:09
Operator : CG/JU
Sample : M2372-09
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CHRT-26516

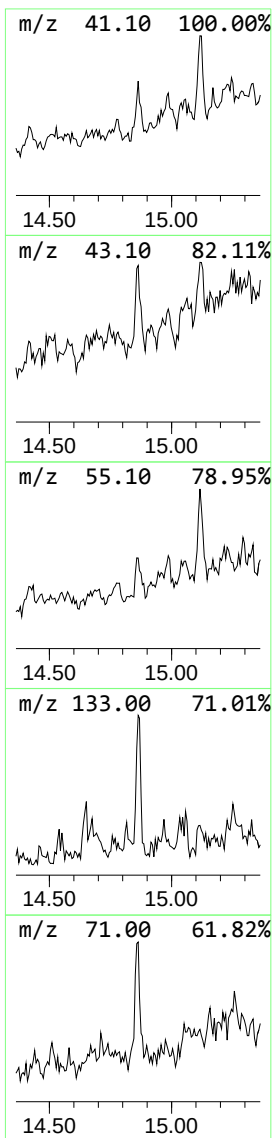
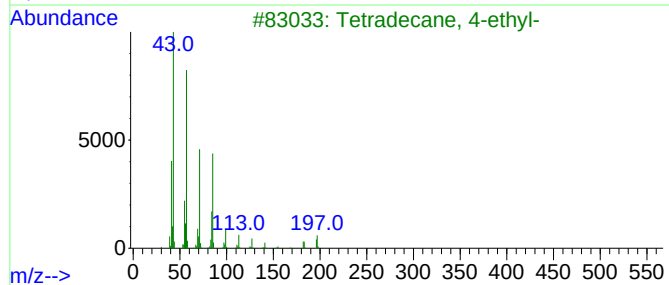
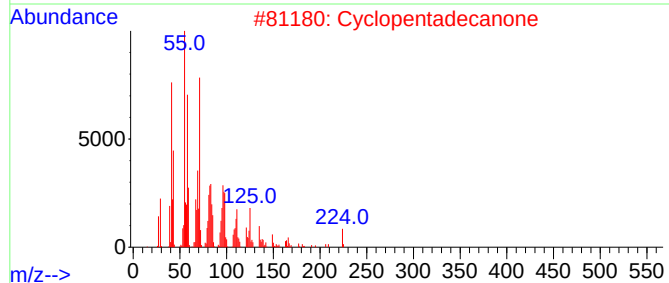
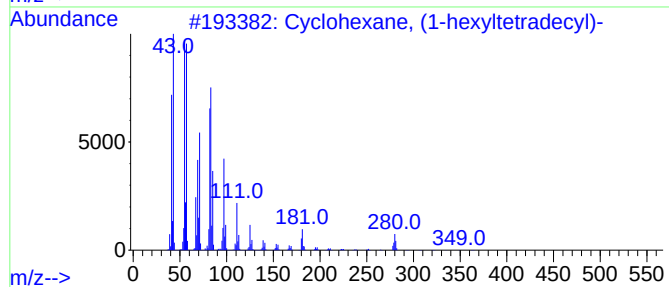
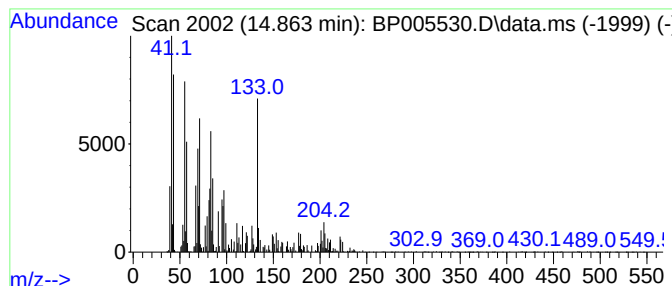
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 unknown14.863 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.863	4.63 ng	62158	Acenaphthene-d10	14.339

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, (1-hexyltetradecyl)-	364	C26H52	004443-60-1	43
2			Cyclopentadecanone	224	C15H28O	000502-72-7	35
3			Tetradecane, 4-ethyl-	226	C16H34	055045-14-2	25
4			Tridecane, 3-ethyl-	212	C15H32	013286-73-2	25
5			1-Iodo-2-methylundecane	296	C12H25I	073105-67-6	18



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051321\
Data File : BP005530.D
Acq On : 13 May 2021 21:09
Operator : CG/JU
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Misc :
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Instrument :
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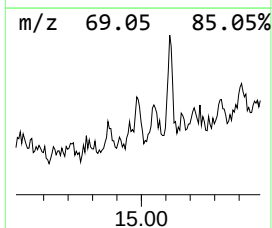
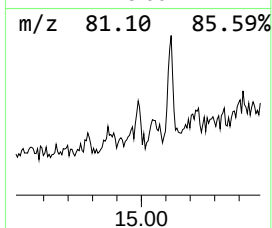
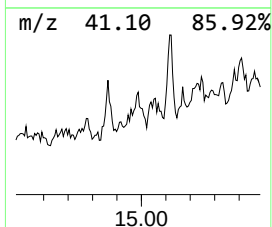
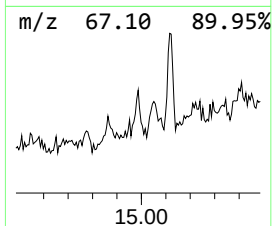
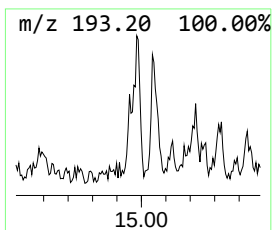
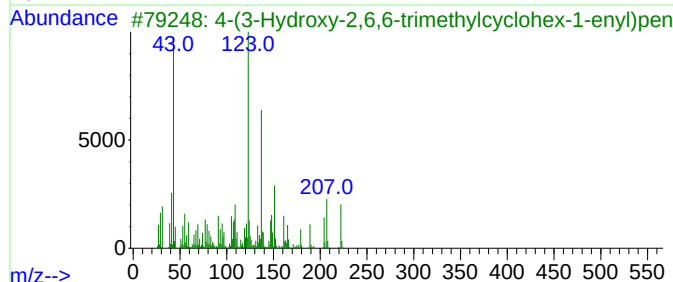
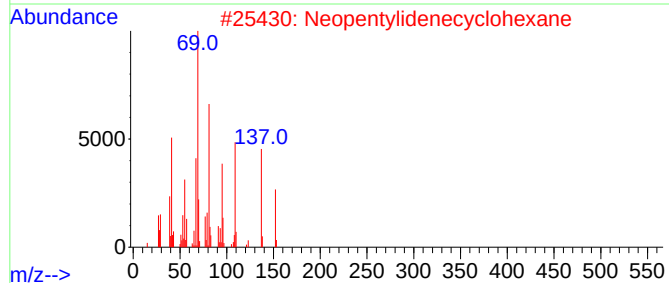
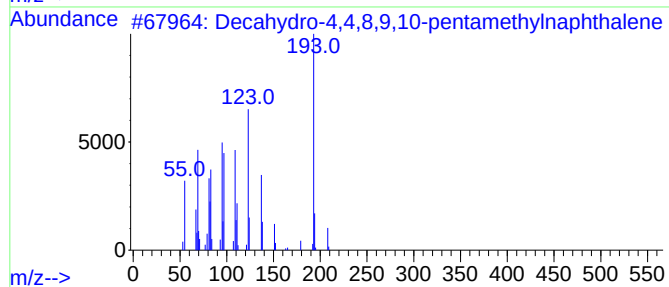
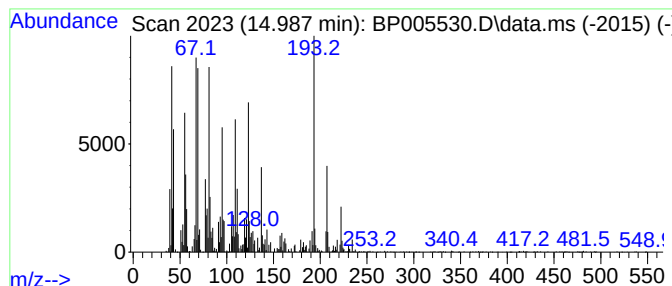
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 unknown14.987 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.987	8.67 ng	116371	Acenaphthene-d10	14.339

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Decahydro-4,4,8,9,10-pentamethyl...	208	C15H28	080655-44-3	52
2			Neopentylidenecyclohexane	152	C11H20	039546-80-0	43
3			4-(3-Hydroxy-2,6,6-trimethylcycl...	222	C14H22O2	097306-57-5	25
4			Spiro[2,4,5,6,7,7a-hexahydro-2-o...	208	C12H16O3	1000197-10-9	22
5			2-Isopropylidene-5-methylhex-4-enal	152	C10H16O	003304-28-7	22



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051321\
Data File : BP005530.D
Acq On : 13 May 2021 21:09
Operator : CG/JU
Sample : M2372-09
Misc :
ALS Vial : 10 Sample Multiplier: 1

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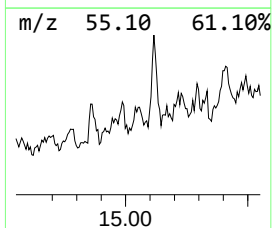
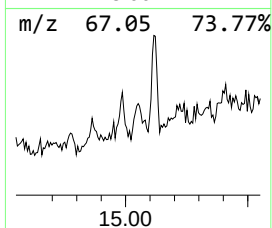
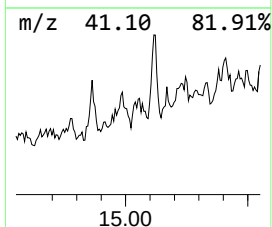
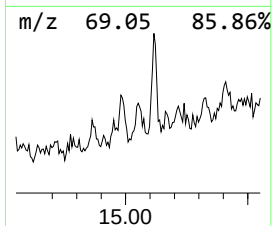
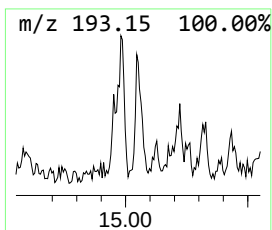
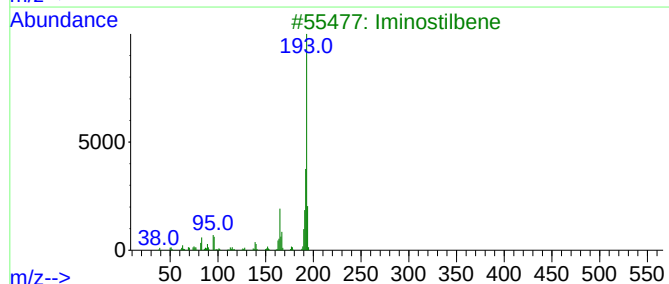
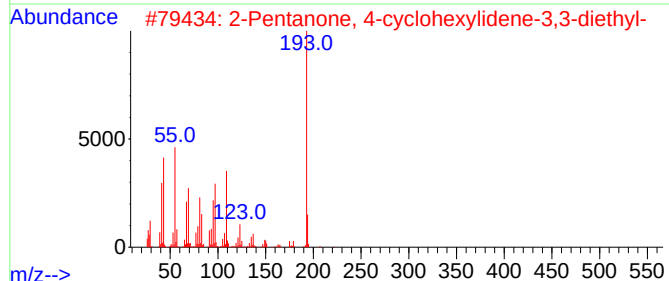
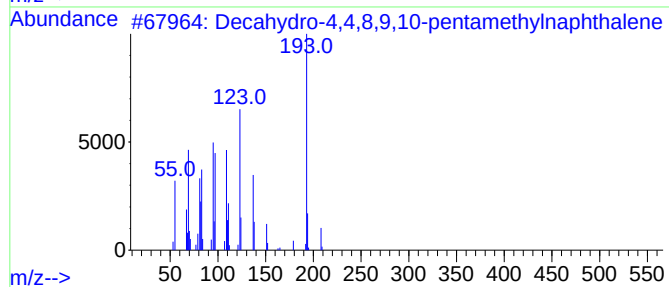
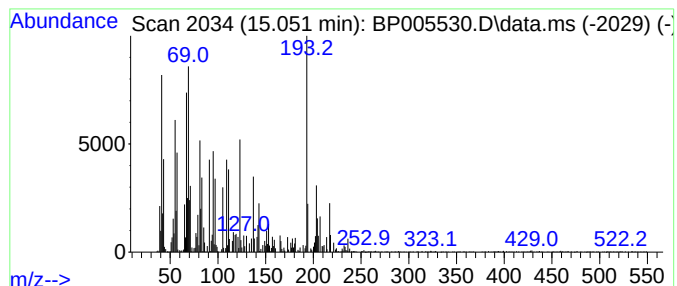
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 unknown15.051 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.051	8.98 ng	120580	Acenaphthene-d10	14.339

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Decahydro-4,4,8,9,10-pentamethyl...	208	C15H28	080655-44-3	38
2			2-Pentanone, 4-cyclohexylidene-3...	222	C15H26O	313253-65-5	27
3			Iminostilbene	193	C14H11N	000256-96-2	14
4			4H-1,3,2-Dioxaborin, 4-ethenyl-2...	208	C12H21B02	074630-04-9	12
5			2(1H)-Naphthalenone, octahydro-4...	194	C13H22O	007056-56-6	11



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051321\
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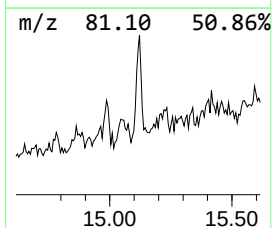
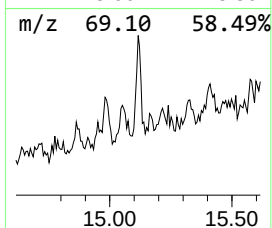
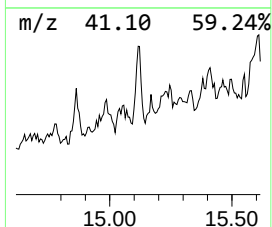
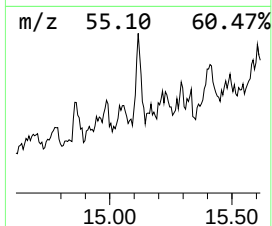
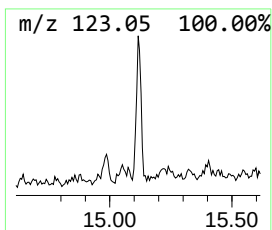
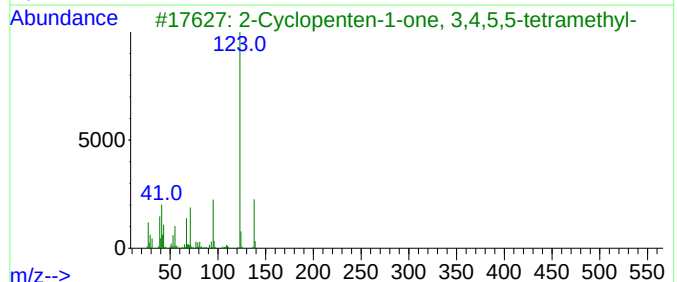
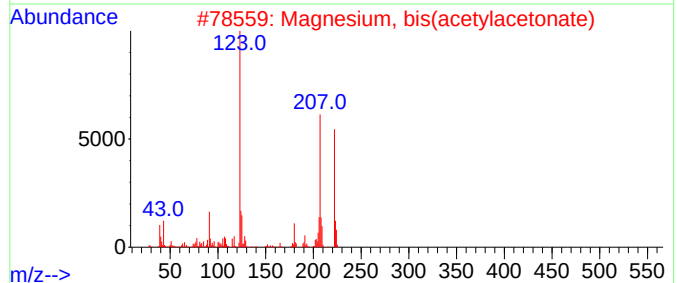
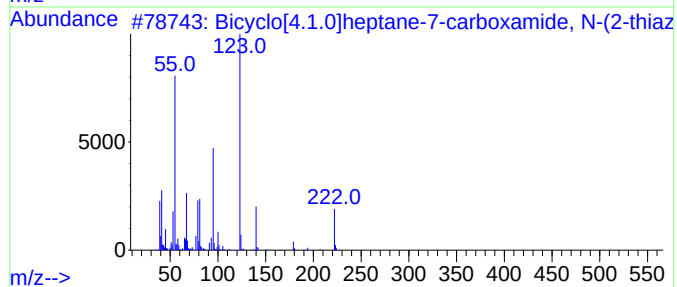
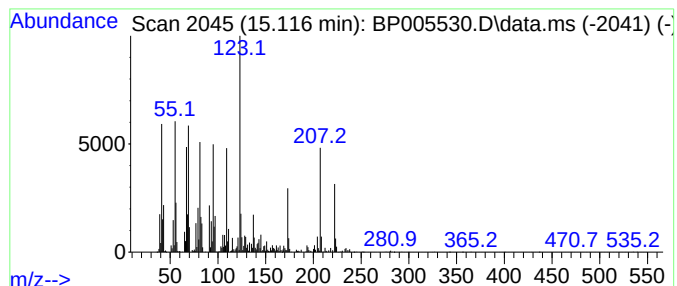
Instrument :
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Bicyclo[4.1.0]heptane-7-car... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD		R.T.	
15.116	15.58 ng	209117	Acenaphthene-d10		14.339	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Bicyclo[4.1.0]heptane-7-carboxam...		222	C11H14N2O5	329912-72-3	55
2	Magnesium, bis(acetylacetonate)		222	C10H14MgO4	068488-07-3	43
3	2-Cyclopenten-1-one, 3,4,5,5-tet...		138	C9H14O	090611-02-2	38
4	2-Pyrimidinamine, 4,6-dimethyl-		123	C6H9N3	000767-15-7	38
5	Bicyclo[2.2.1]heptane, 1,3,3-tri...		138	C10H18	006248-88-0	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051321\
 Data File : BP005530.D
 Acq On : 13 May 2021 21:09
 Operator : CG/JU
 Sample : M2372-09
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
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 ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP050521.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
Decahydro-4,4,8...	13.098	3.8	ng	50633	3	14.339	268507	20.0
1H-Indene, octa...	13.746	5.8	ng	77829	3	14.339	268507	20.0
unknown13.863	13.863	4.6	ng	62349	3	14.339	268507	20.0
unknown14.063	14.063	4.8	ng	64472	3	14.339	268507	20.0
unknown14.116	14.116	2.5	ng	33536	3	14.339	268507	20.0
2-Pentanone, 4-...	14.151	10.6	ng	142706	3	14.339	268507	20.0
unknown14.234	14.234	4.1	ng	54792	3	14.339	268507	20.0
unknown14.287	14.287	6.3	ng	84143	3	14.339	268507	20.0
unknown14.863	14.863	4.6	ng	62158	3	14.339	268507	20.0
unknown14.987	14.987	8.7	ng	116371	3	14.339	268507	20.0
unknown15.051	15.051	9.0	ng	120580	3	14.339	268507	20.0
Bicyclo[4.1.0]h...	15.116	15.6	ng	209117	3	14.339	268507	20.0