

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041221\
 Data File : BP005274.D
 Acq On : 12 Apr 2021 19:41
 Operator : CG/JU
 Sample : M1997-02
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 RBR-58-56

Integration Parameters: rteint.p

Integrator: RTE

Smothing : 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-BP040921.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BP005274.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	367	373	382	rVB	853137	1242037	67.81%	9.094%
2	6.199	525	529	540	rVB2	43493	79282	4.33%	0.581%
3	6.863	636	642	651	rVB	736312	1139963	62.24%	8.347%
4	6.952	652	657	659	rBV	168145	262483	14.33%	1.922%
5	7.675	775	780	789	rVB	166179	272577	14.88%	1.996%
6	8.340	888	893	900	rVB3	15618	22566	1.23%	0.165%
7	8.840	971	978	986	rBV	412394	681263	37.20%	4.988%
8	8.910	986	990	996	rVB2	17344	27536	1.50%	0.202%
9	9.169	1028	1034	1043	rVB	55783	89736	4.90%	0.657%
10	10.463	1247	1254	1262	rBV	180343	301368	16.45%	2.207%
11	12.951	1667	1677	1684	rBV	863753	1237672	67.58%	9.063%
12	13.104	1698	1703	1707	rBV6	13033	23996	1.31%	0.176%
13	13.598	1784	1787	1791	rVB5	17527	21451	1.17%	0.157%
14	13.745	1810	1812	1816	rVB4	20551	22434	1.22%	0.164%
15	13.798	1818	1821	1823	rBV3	24844	35153	1.92%	0.257%
16	13.863	1829	1832	1835	rBV4	31444	41828	2.28%	0.306%
17	14.163	1879	1883	1890	rVB5	84817	150945	8.24%	1.105%
18	14.239	1893	1896	1901	rVB3	90568	126702	6.92%	0.928%
19	14.339	1909	1913	1920	rVB	301511	456213	24.91%	3.341%
20	15.198	2056	2059	2063	rBV	143733	188418	10.29%	1.380%
21	15.845	2164	2169	2173	rBV	687954	981822	53.61%	7.189%
22	16.069	2204	2207	2215	rVB	212505	344963	18.83%	2.526%
23	16.869	2340	2343	2347	rVB	163950	170038	9.28%	1.245%
24	17.104	2379	2383	2388	rBV	382482	543905	29.70%	3.983%
25	18.292	2581	2585	2592	rBV2	648754	915167	49.97%	6.701%
26	19.680	2817	2821	2827	rVB	1599803	1831529	100.00%	13.411%
27	20.310	2924	2928	2932	rBV	356111	471598	25.75%	3.453%
28	20.916	3028	3031	3039	rVB3	255853	451639	24.66%	3.307%
29	20.986	3040	3043	3052	rVB	203204	255900	13.97%	1.874%
30	21.204	3075	3080	3084	rBV	444094	577551	31.53%	4.229%
31	21.839	3185	3188	3198	rVB	186635	250158	13.66%	1.832%
32	23.474	3461	3466	3472	rBV	248300	439123	23.98%	3.215%

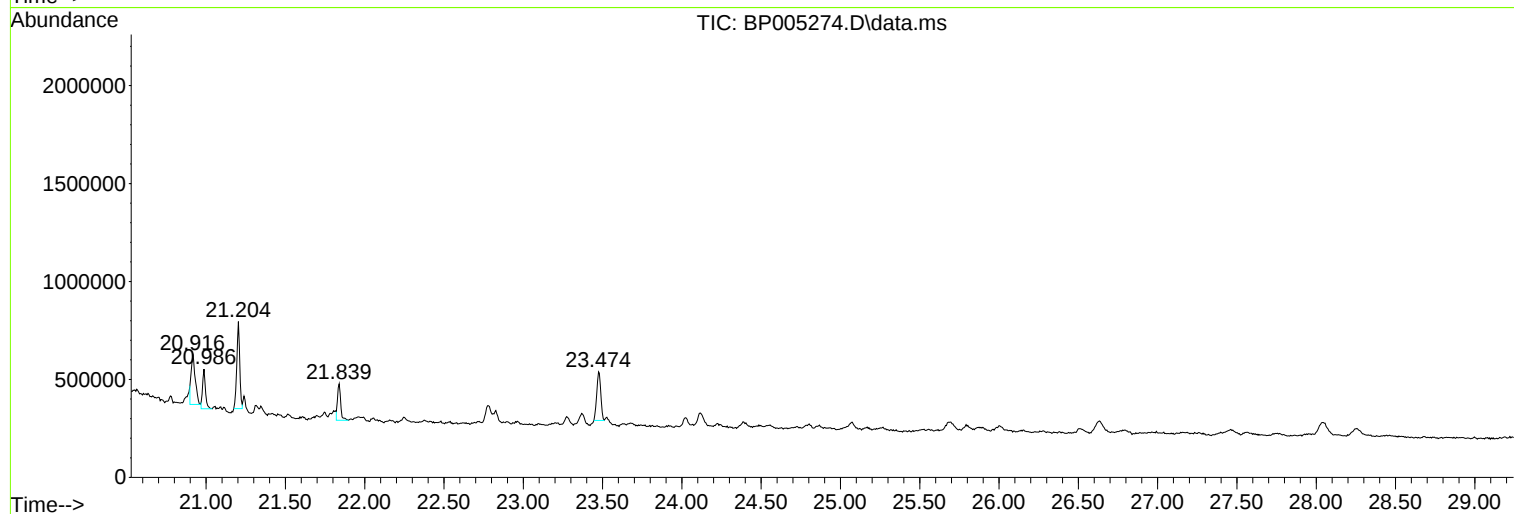
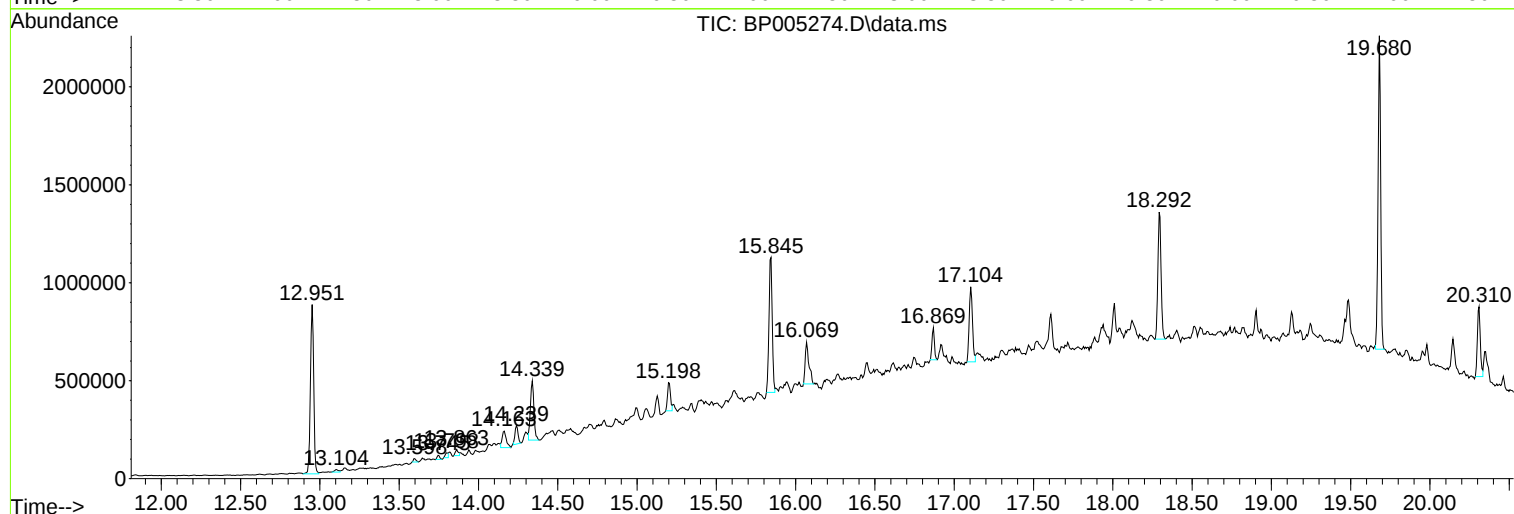
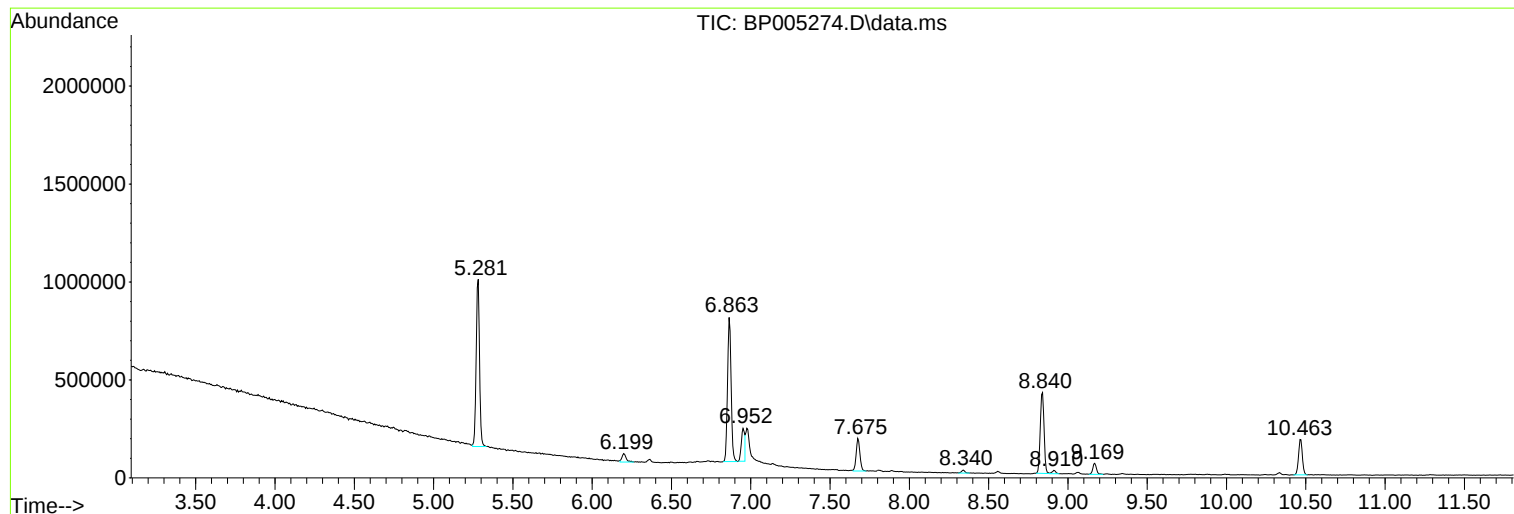
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP040921.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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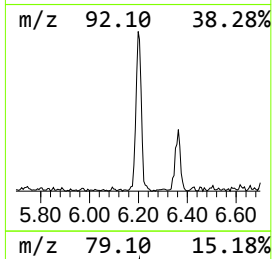
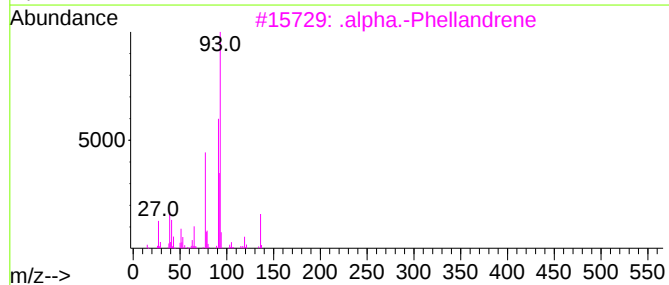
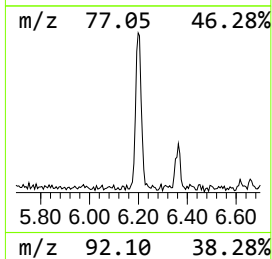
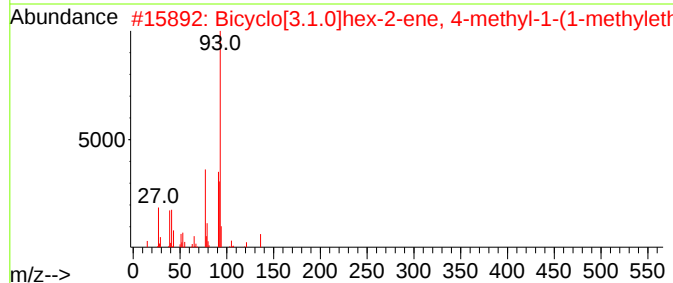
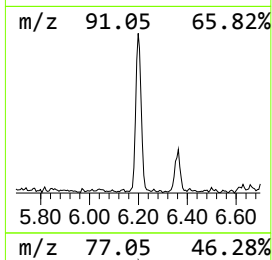
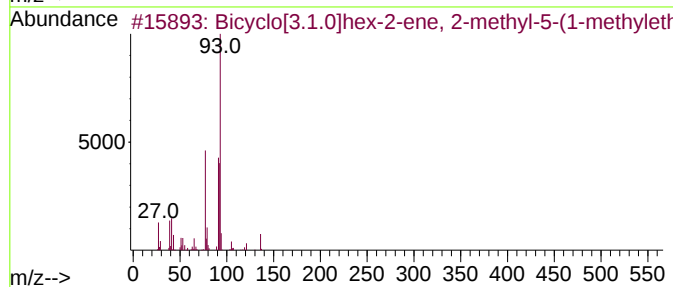
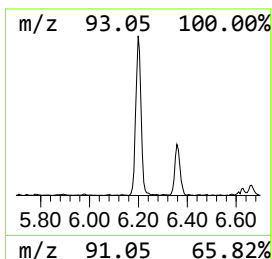
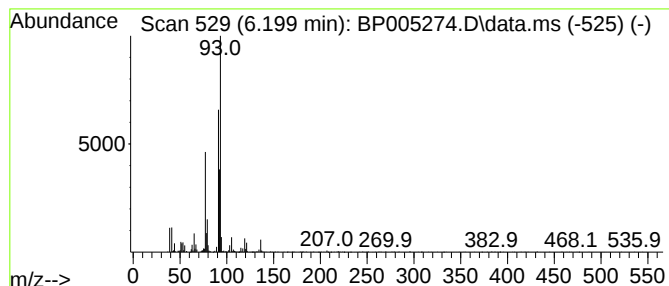
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TIC Integration Parameters: LSCINT.P

Peak Number 1 Bicyclo[3.1.0]hex-2-ene, 2-... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.199	5.82 ng	79282	1,4-Dichlorobenzene-d4	7.675

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Bicyclo[3.1.0]hex-2-ene, 2-methy...	136	C10H16	002867-05-2	91
2			Bicyclo[3.1.0]hex-2-ene, 4-methy...	136	C10H16	028634-89-1	90
3			.alpha.-Phellandrene	136	C10H16	000099-83-2	87
4			1,3-Cyclohexadiene, 5,6-dimethyl-	108	C8H12	005715-27-5	53
5			Pyridine, 2-propyl-	121	C8H11N	000622-39-9	43



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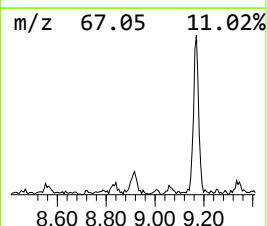
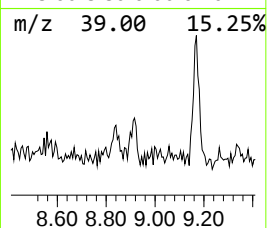
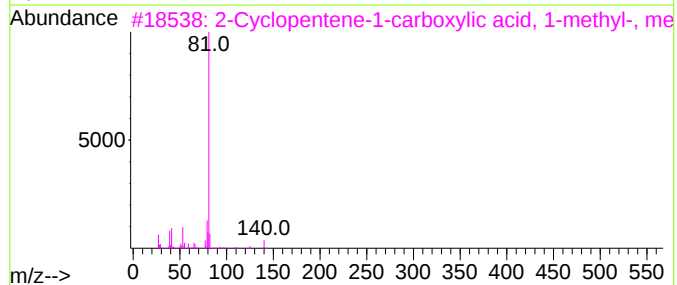
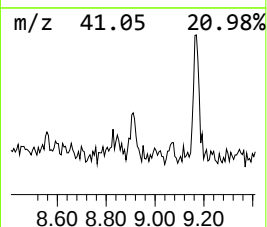
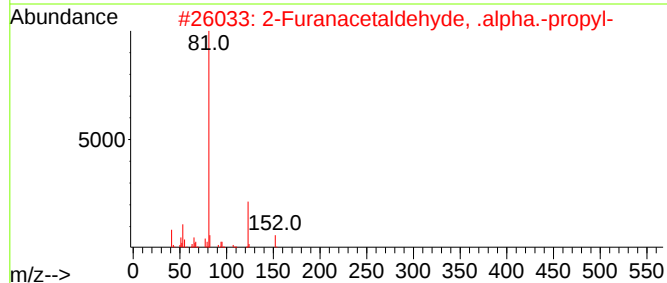
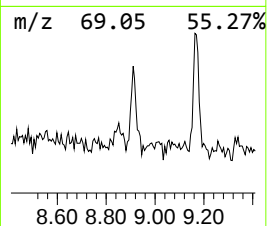
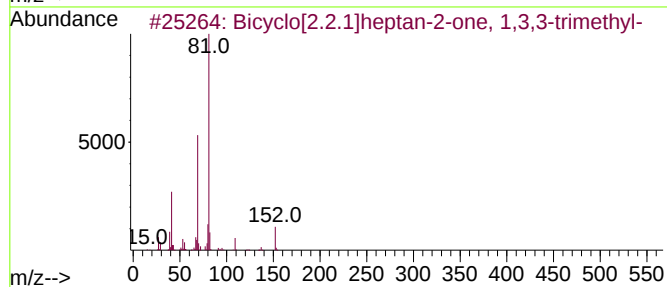
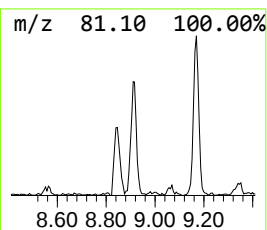
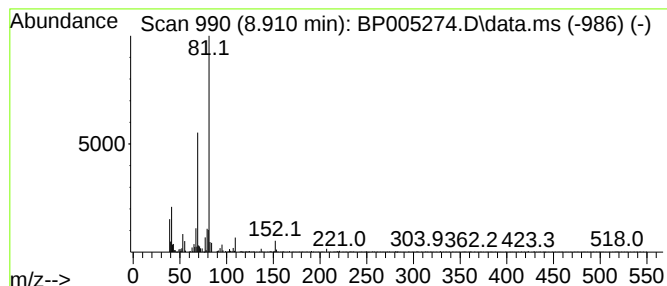
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 2 Bicyclo[2.2.1]heptan-2-one,... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.910	2.02 ng	27536	1,4-Dichlorobenzene-d4	7.675

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Bicyclo[2.2.1]heptan-2-one, 1,3,...	152	C10H16O	001195-79-5	80
2			2-Furanacetaldehyde, .alpha.-pro...	152	C9H12O2	031681-26-2	64
3			2-Cyclopentene-1-carboxylic acid...	140	C8H12O2	068317-73-7	64
4			Bicyclo[2.1.0]pentane, 1,4-dimet...	96	C7H12	017065-18-8	53
5			2,4-Dimethyl 1,4-pentadiene	96	C7H12	004161-65-3	53



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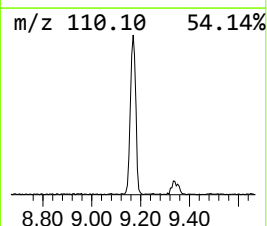
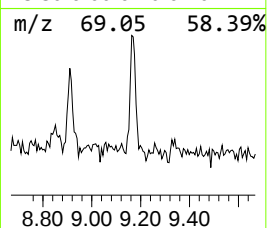
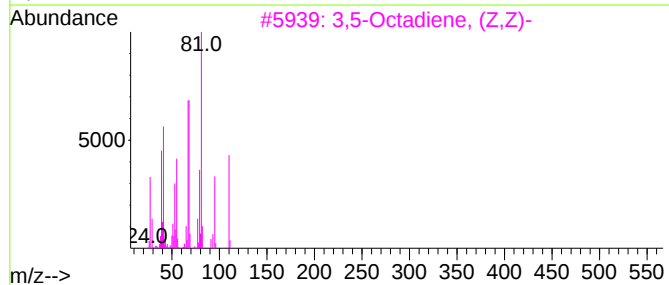
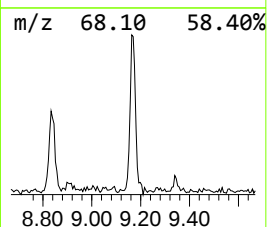
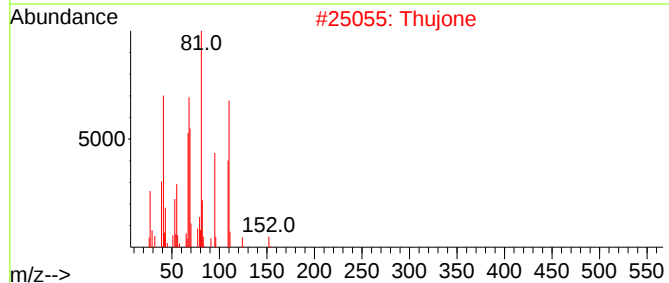
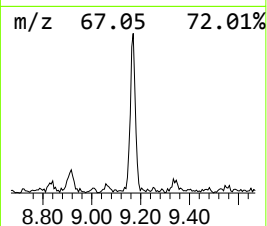
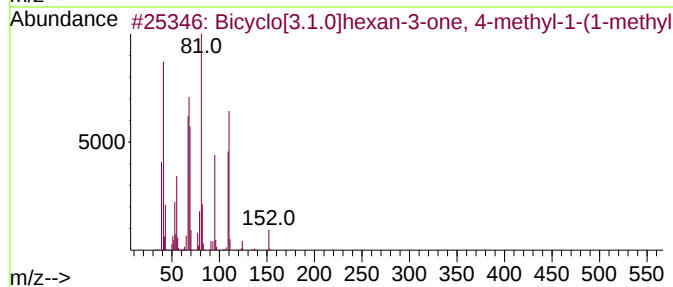
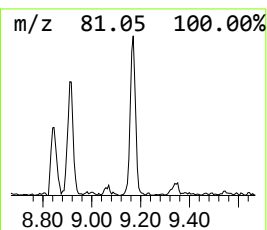
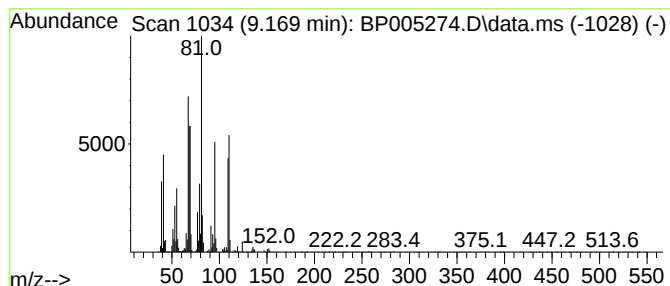
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 3 Bicyclo[3.1.0]hexan-3-one, ... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.169	5.96 ng	89736	Naphthalene-d8	10.469

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Bicyclo[3.1.0]hexan-3-one, 4-met...	152	C10H16O	001125-12-8	91
2			Thujone	152	C10H16O	000546-80-5	90
3			3,5-Octadiene, (Z,Z)-	110	C8H14	007348-80-3	72
4			Methyl ethyl cyclopentene	110	C8H14	019780-56-4	70
5			Cyclohexene, 1,2-dimethyl-	110	C8H14	001674-10-8	62



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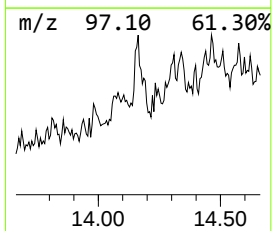
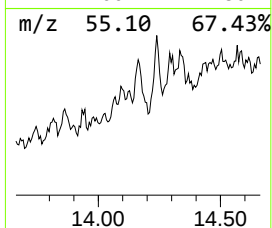
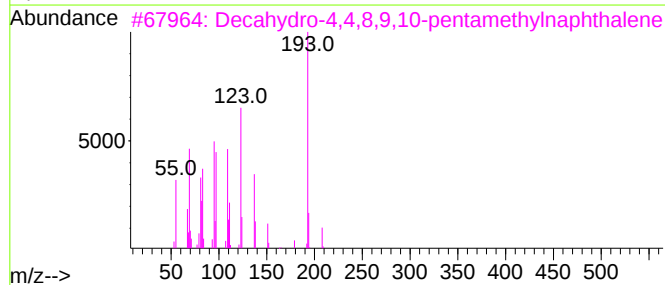
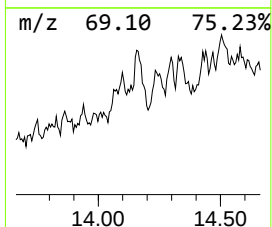
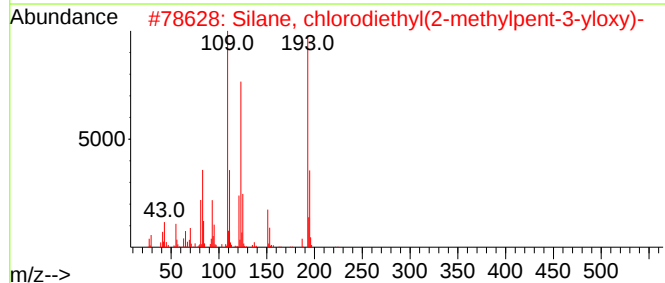
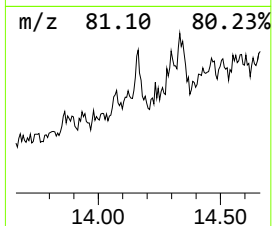
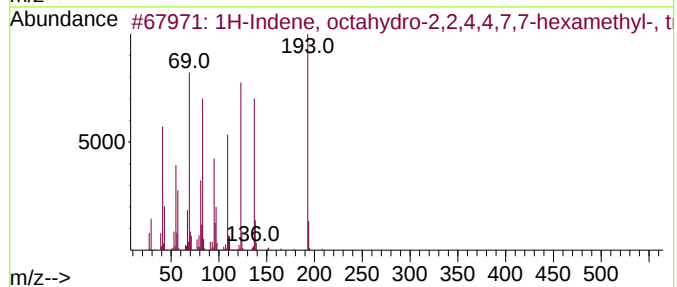
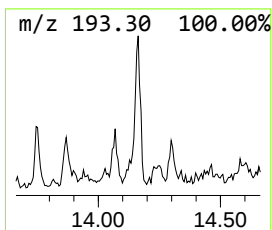
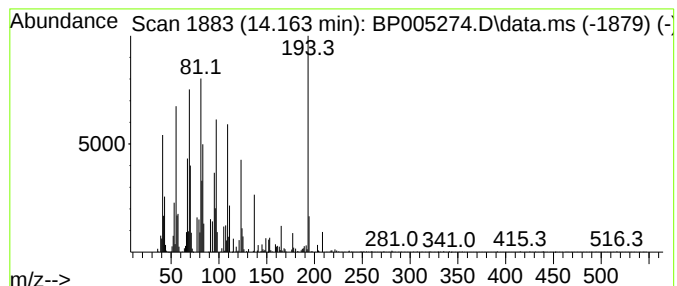
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Peak Number 4 unknown14.163 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.163	6.62 ng	150945	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	38
2			Silane, chlorodiethyl(2-methylpe...	222	C10H23ClOSi	1000363-79-1	22
3			Decahydro-4,4,8,9,10-pentamethyl...	208	C15H28	080655-44-3	22
4			1,22-Docosanediol	342	C22H46O2	022513-81-1	16
5			3,7-Dimethylocta-2,6-dien-1-ol, ...	262	C16H22OS	1000196-46-7	12



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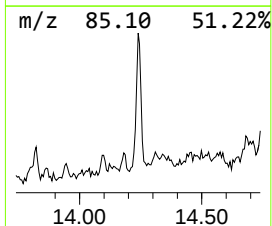
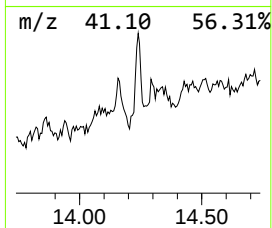
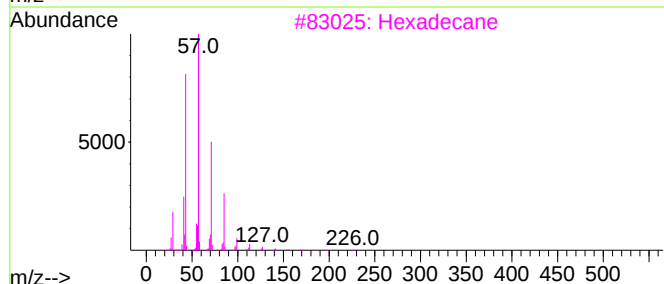
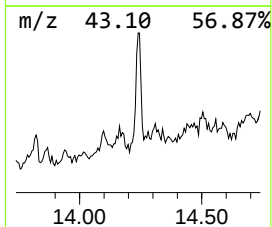
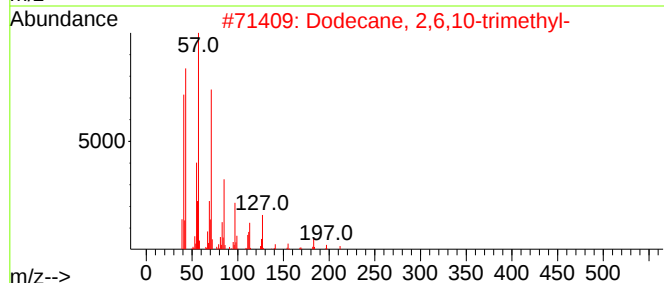
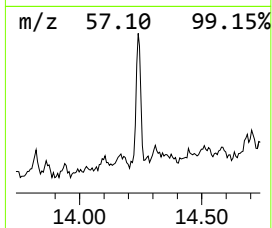
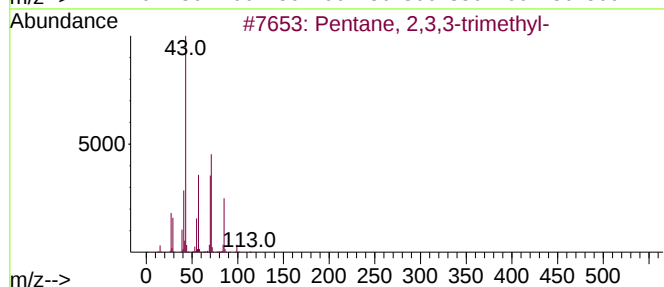
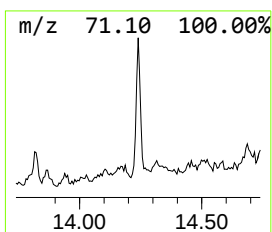
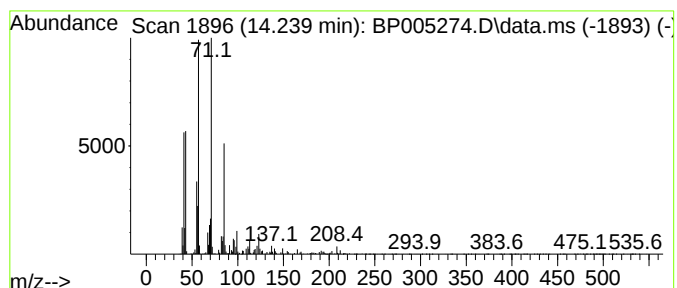
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Peak Number 5 Pentane, 2,3,3-trimethyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.239	5.55 ng	126702	Acenaphthene-d10	14.339

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentane, 2,3,3-trimethyl-	114	C8H18	000560-21-4	59
2			Dodecane, 2,6,10-trimethyl-	212	C15H32	003891-98-3	59
3			Hexadecane	226	C16H34	000544-76-3	59
4			Sulfurous acid, butyl decyl ester	278	C14H30O3S	1000309-17-7	47
5			Dodecane	170	C12H26	000112-40-3	47



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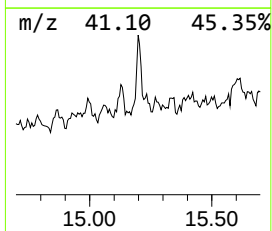
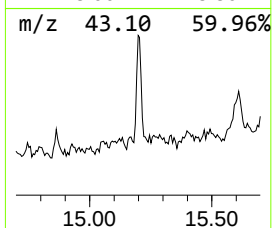
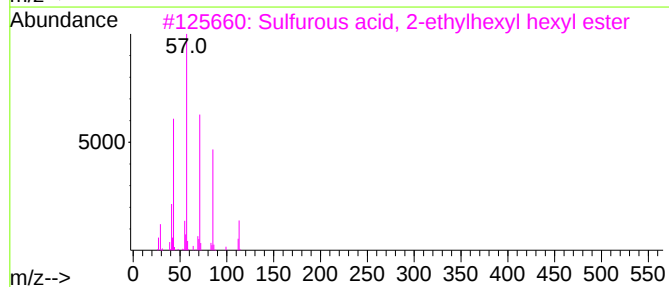
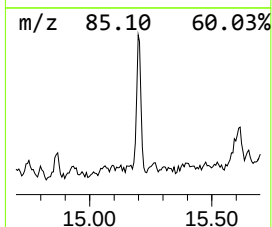
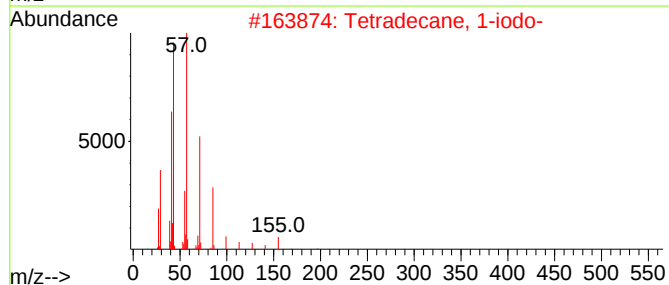
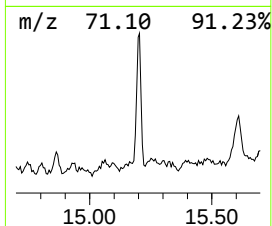
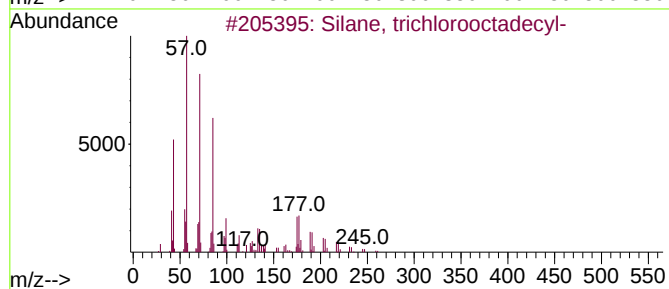
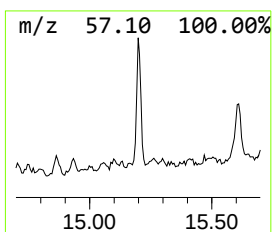
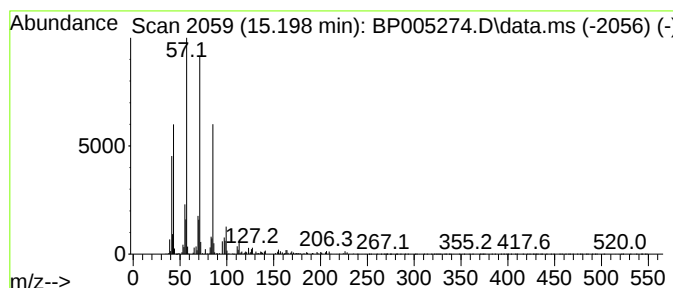
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Silane, trichlorooctadecyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.198	8.26 ng	188418	Acenaphthene-d10	14.339

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Silane, trichlorooctadecyl-	386	C18H37Cl3Si	000112-04-9	72
2		Tetradecane, 1-iodo-	324	C14H29I	019218-94-1	72
3		Sulfurous acid, 2-ethylhexyl hex...	278	C14H30O3S	1000309-20-2	64
4		Sulfurous acid, 2-ethylhexyl tri...	376	C21H44O3S	1000309-19-6	64
5		Sulfurous acid, pentyl undecyl e...	306	C16H34O3S	1000309-14-4	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041221\
Data File : BP005274.D
Acq On : 12 Apr 2021 19:41
Operator : CG/JU
Sample : M1997-02
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
RBR-58-56

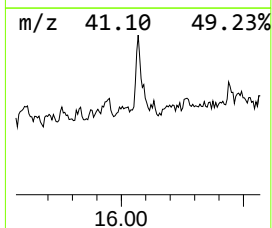
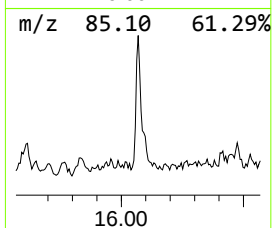
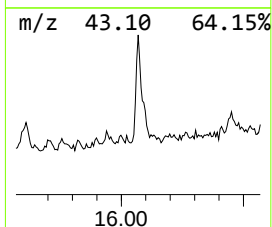
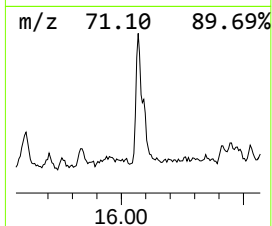
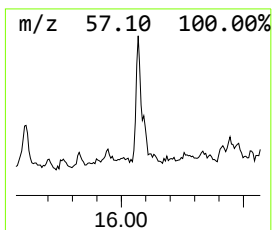
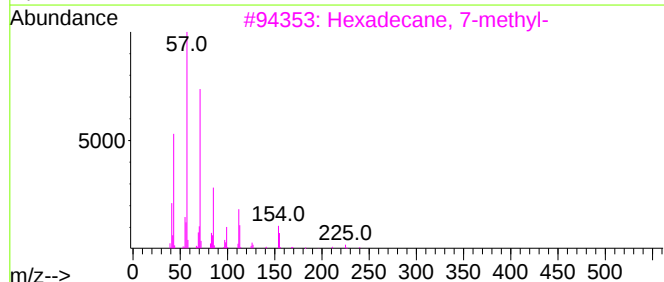
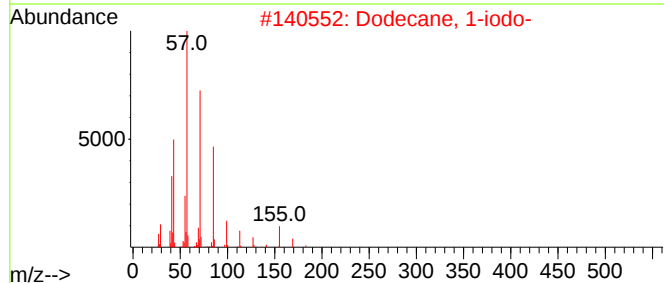
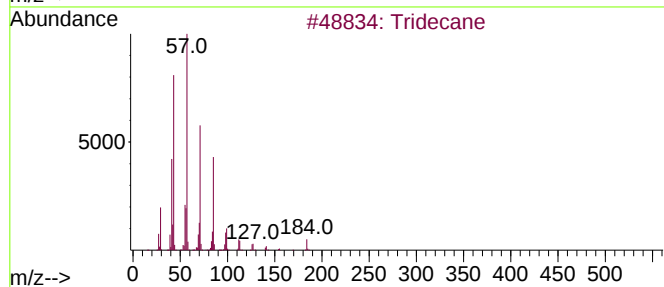
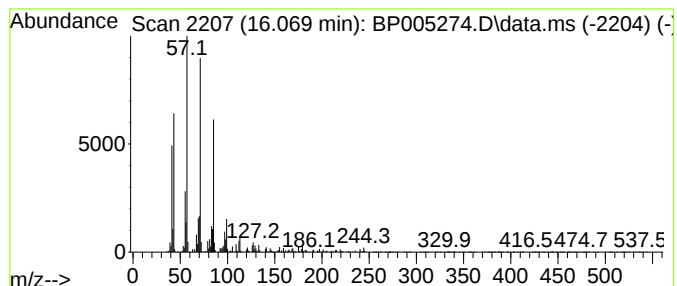
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Tridecane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.069	12.68 ng	344963	Phenanthrene-d10	17.104

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tridecane	184	C13H28	000629-50-5	83
2			Dodecane, 1-iodo-	296	C12H25I	004292-19-7	80
3			Hexadecane, 7-methyl-	240	C17H36	026730-20-1	72
4			Dodecane, 4,6-dimethyl-	198	C14H30	061141-72-8	70
5			Tetradecane, 2,6,10-trimethyl-	240	C17H36	014905-56-7	68



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041221\
Data File : BP005274.D
Acq On : 12 Apr 2021 19:41
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Sample : M1997-02
Misc :
ALS Vial : 15 Sample Multiplier: 1

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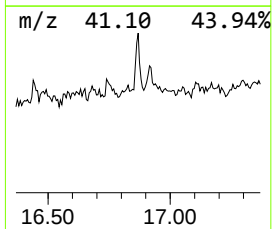
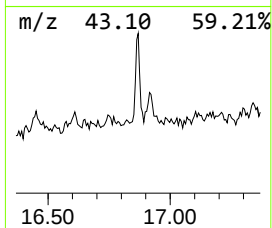
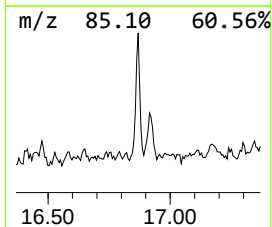
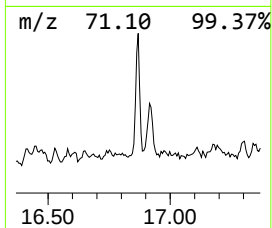
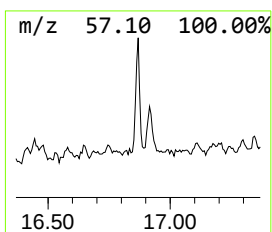
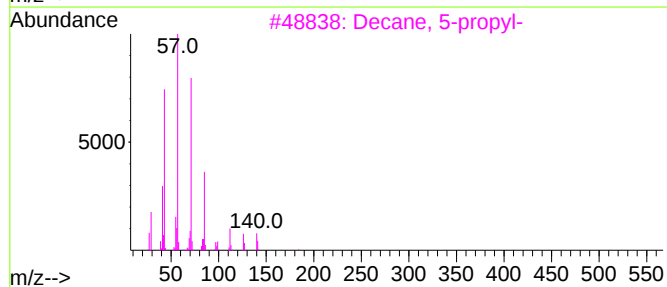
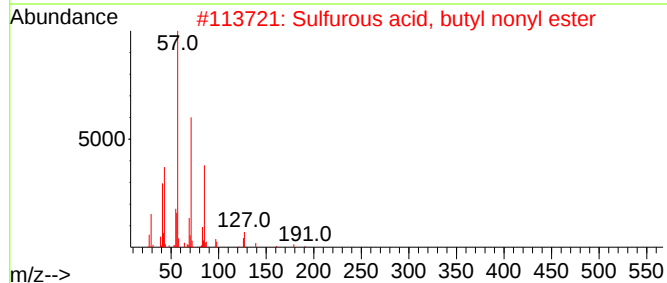
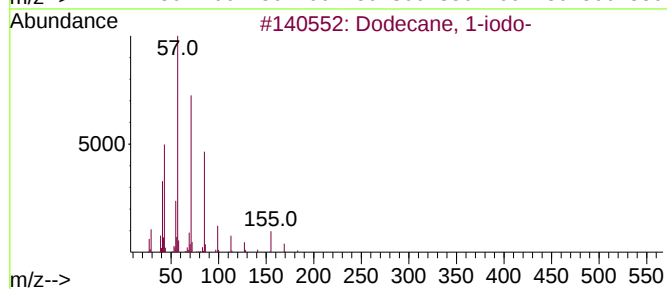
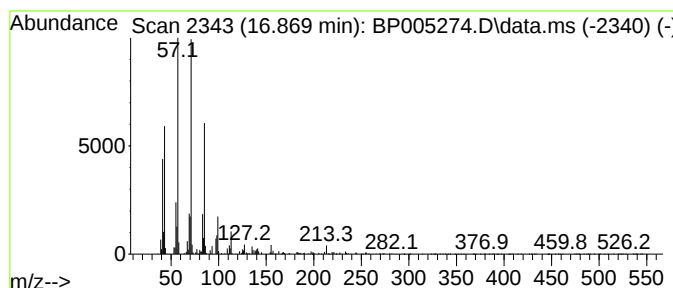
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Dodecane, 1-iodo- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.869	6.25 ng	170038	Phenanthrene-d10	17.104

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dodecane, 1-iodo-	296	C12H25I	004292-19-7	86
2			Sulfurous acid, butyl nonyl ester	264	C13H28O3S	1000309-17-6	72
3			Decane, 5-propyl-	184	C13H28	017312-62-8	58
4			Sulfurous acid, hexyl pentyl ester	236	C11H24O3S	1000309-14-1	58
5			Sulfurous acid, pentyl tridecyl ...	334	C18H38O3S	1000309-14-6	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041221\
Data File : BP005274.D
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Operator : CG/JU
Sample : M1997-02
Misc :
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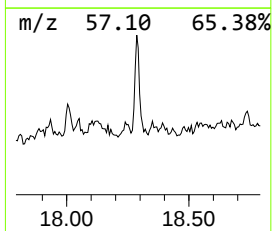
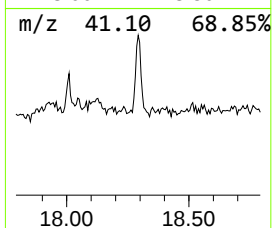
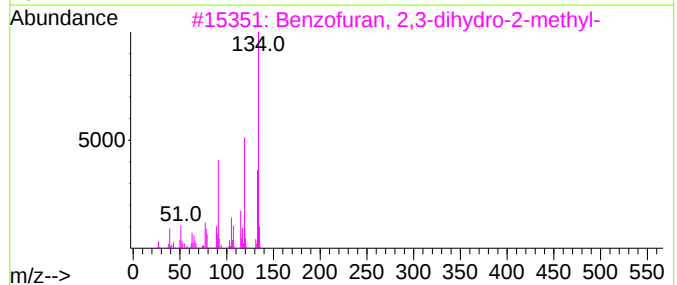
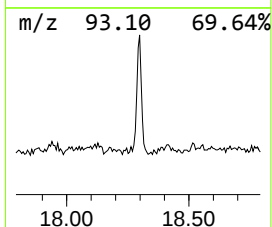
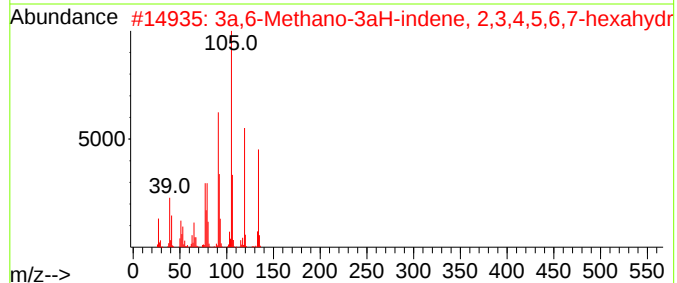
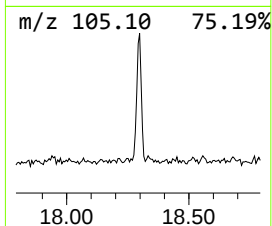
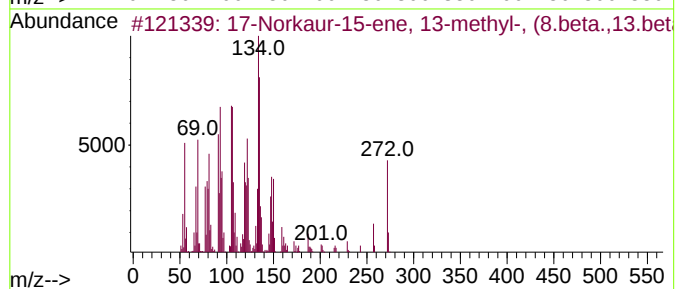
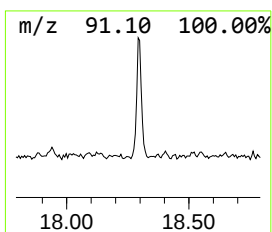
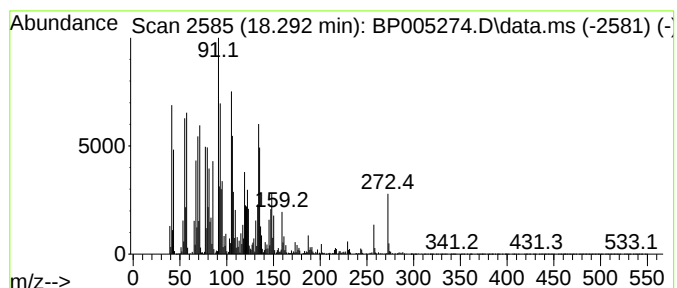
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 17-Norkaur-15-ene, 13-methy... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD			R.T.
18.292	33.65 ng	915167	Phenanthrene-d10			17.104
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	17-Norkaur-15-ene, 13-methyl-, (...)		272	C20H32	003564-54-3	96
2	3a,6-Methano-3aH-indene, 2,3,4,5...		134	C10H14	098640-10-9	25
3	Benzofuran, 2,3-dihydro-2-methyl-		134	C9H10O	001746-11-8	15
4	Benzene, 1-methyl-3-propyl-		134	C10H14	001074-43-7	14
5	Succinic acid, 2-adamantyl ethyl...		280	C16H24O4	1000329-93-1	14



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041221\
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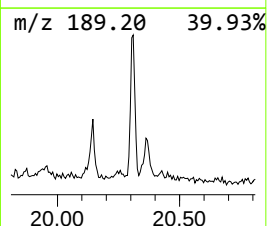
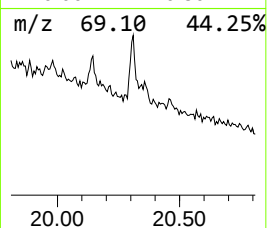
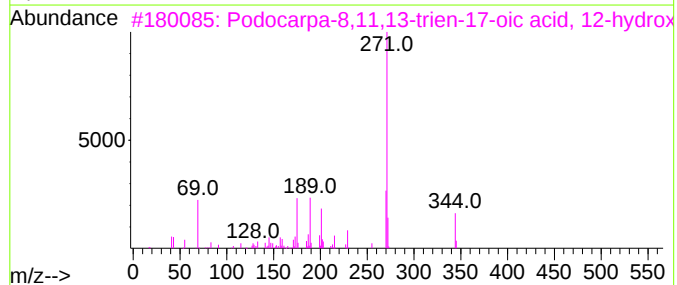
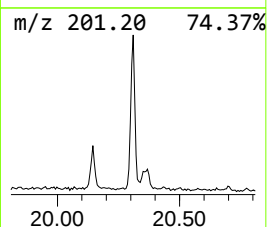
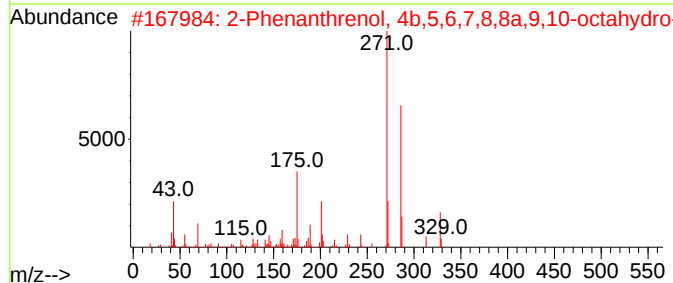
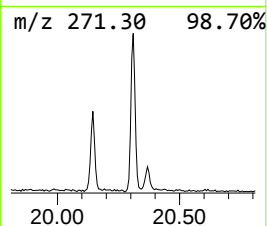
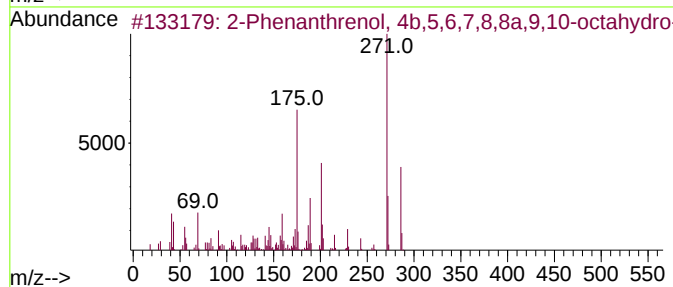
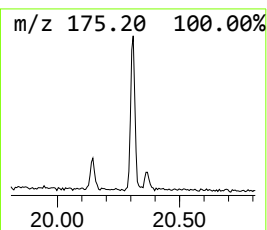
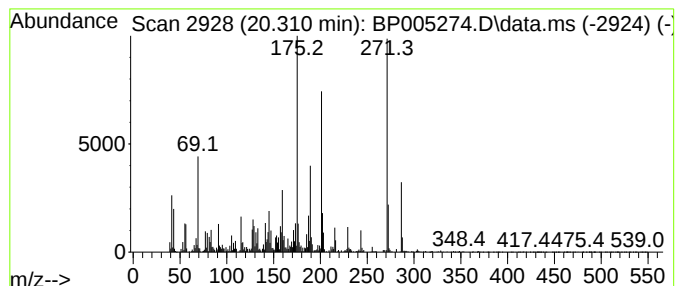
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 2-Phenanthrenol, 4b,5,6,7,8... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD		R.T.	
20.310	16.33 ng	471598	Chrysene-d12		21.204	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	2-Phenanthrenol, 4b,5,6,7,8,8a,9...		286	C20H30O	000511-15-9	83
2	2-Phenanthrenol, 4b,5,6,7,8,8a,9...		328	C22H32O2	015340-82-6	81
3	Podocarpa-8,11,13-trien-17-oic a...		344	C22H32O3	024067-43-4	46
4	Pyrrolidine, 1-[5-(1,3-benzodiox...		271	C16H17NO3	025924-78-1	35
5	n-Hexyl-6-methoxy-4-methyl-8-qui...		272	C17H24N2O	083547-16-4	18



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041221\
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Acq On : 12 Apr 2021 19:41
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Misc :
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ClientSampleId :
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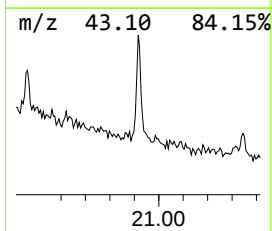
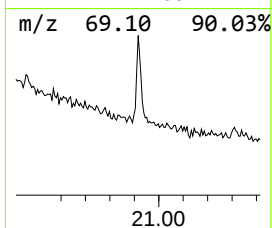
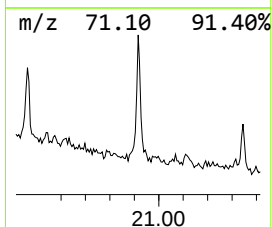
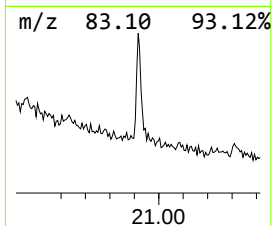
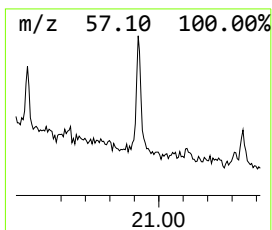
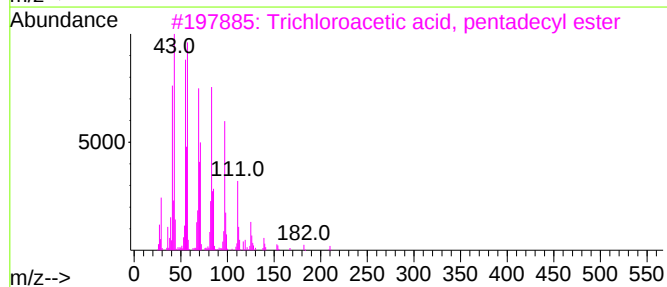
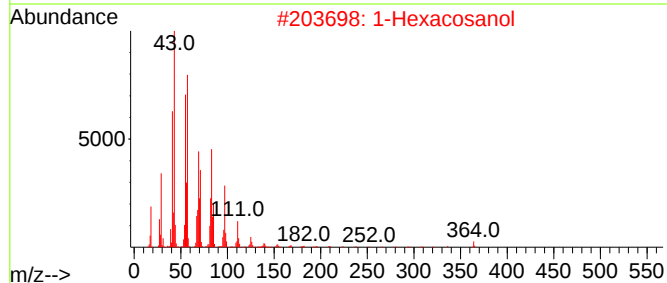
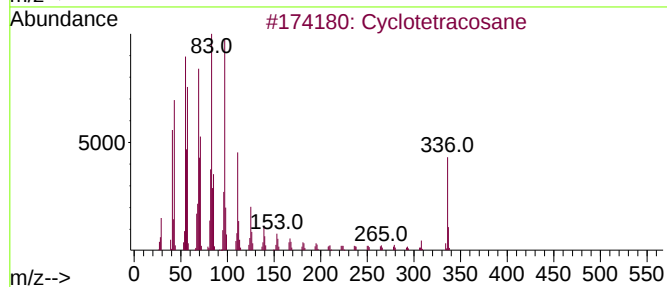
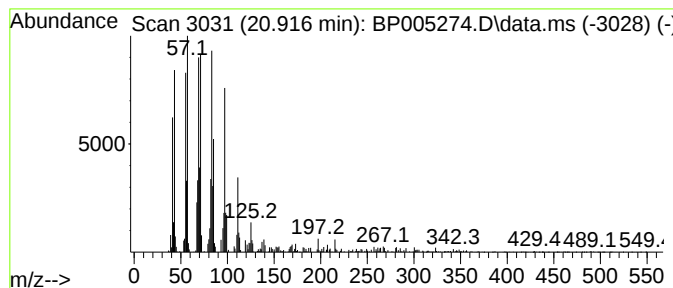
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Cyclotetracosane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.916	15.64 ng	451639	Chrysene-d12	21.204

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclotetracosane	336	C24H48	000297-03-0	89
2			1-Hexacosanol	382	C26H54O	000506-52-5	87
3			Trichloroacetic acid, pentadecyl...	372	C17H31Cl3O2	074339-53-0	83
4			Behenic alcohol	326	C22H46O	000661-19-8	81
5			5-Eicosene, (E)-	280	C20H40	074685-30-6	78



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ClientSampleId :
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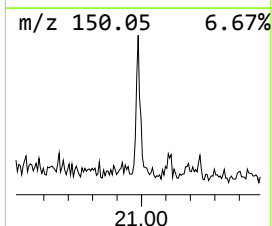
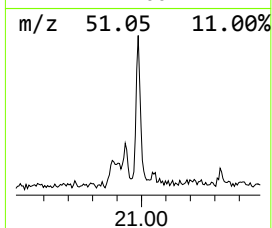
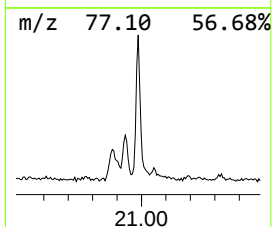
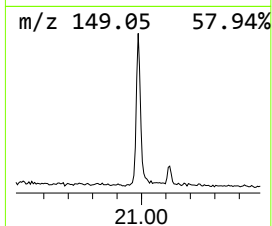
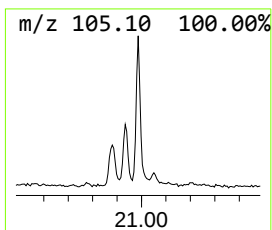
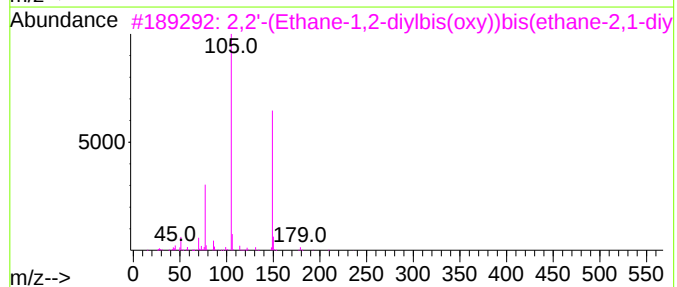
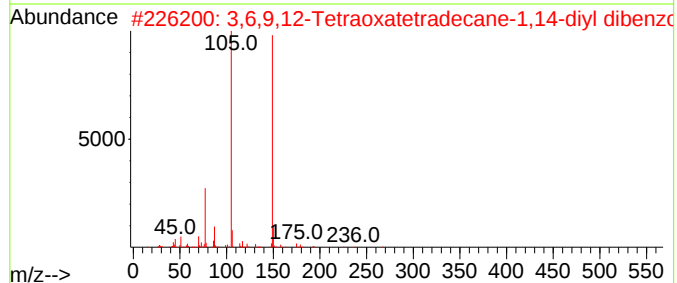
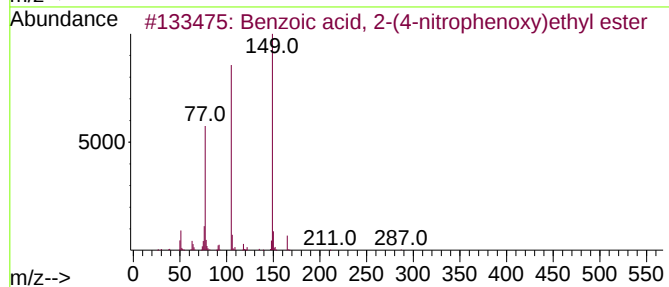
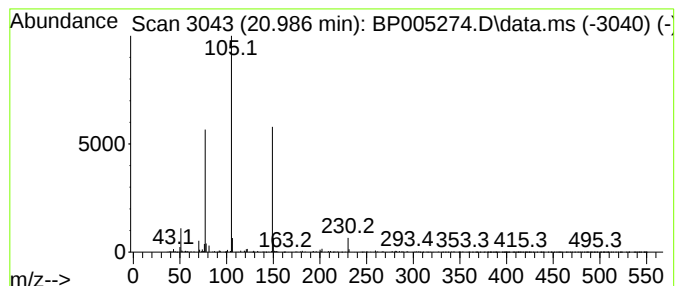
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP040921.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Benzoic acid, 2-(4-nitrophe... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.986	8.86 ng	255900	Chrysene-d12	21.204

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzoic acid, 2-(4-nitrophenoxy)...	287	C15H13NO5	1000368-92-7	72
2			3,6,9,12-Tetraoxatetradecane-1,1...	446	C24H30O8	1000366-97-0	64
3			2,2'-(Ethane-1,2-diylbis(oxy))bi...	358	C20H22O6	1000366-99-4	59
4			Diethylene glycol dibenzoate	314	C18H18O5	000120-55-8	50
5			2-(2-Carboxyvinyl)pyridine, trans	149	C8H7NO2	007340-22-9	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041221\
Data File : BP005274.D
Acq On : 12 Apr 2021 19:41
Operator : CG/JU
Sample : M1997-02
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
RBR-58-56

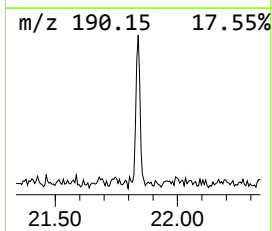
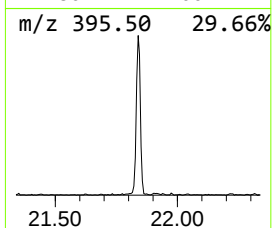
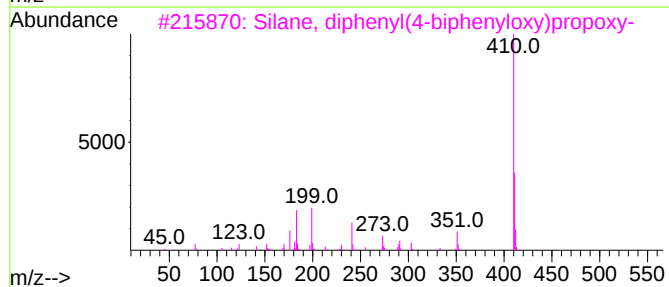
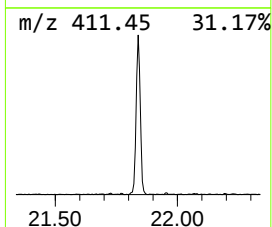
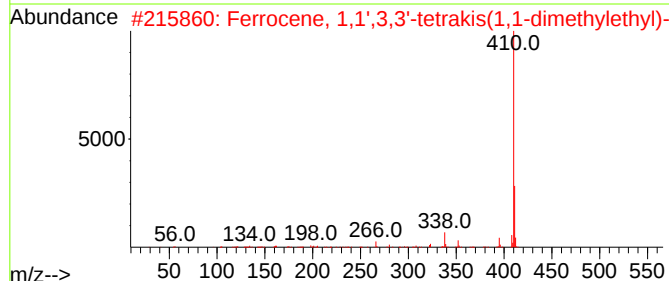
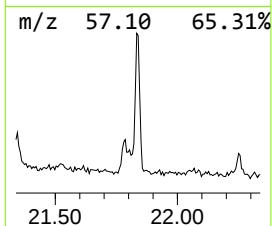
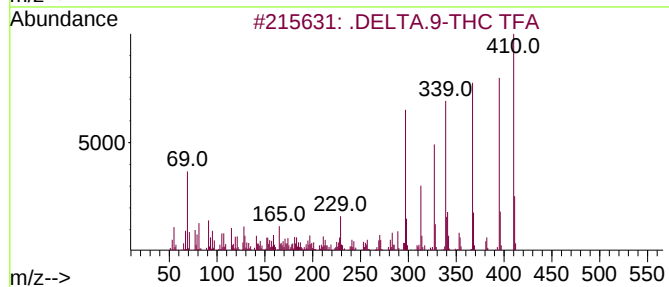
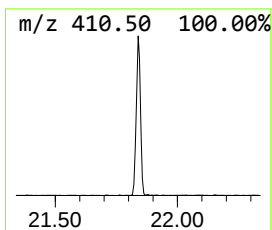
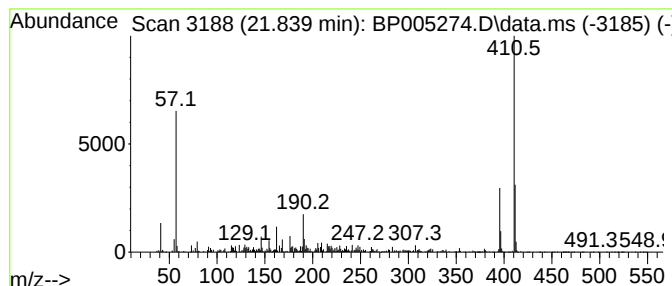
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP040921.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 .DELTA.9-THC TFA Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.839	8.66 ng	250158	Chrysene-d12	21.204

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			.DELTA.9-THC TFA	410	C23H29F3O3	086702-79-6	64
2			Ferrocene, 1,1',3,3'-tetrakis(1,...	410	C26H42Fe	001317-17-5	52
3			Silane, diphenyl(4-biphenyloxy)p...	410	C27H26O2Si	1000367-49-5	28
4			Apomorphine, bis(trimethylsilyl)-	411	C23H33NO2Si2	074841-68-2	9
5			9-Chloro-4,5-dihydro-4-[3,4,5-tr...	410	C22H19ClN2O4	1000212-48-1	9



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041221\
 Data File : BP005274.D
 Acq On : 12 Apr 2021 19:41
 Operator : CG/JU
 Sample : M1997-02
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 RBR-58-56

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP040921.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
Bicyclo[3.1.0]h...	6.199	5.8	ng	79282	1	7.675	272577	20.0
Bicyclo[2.2.1]h...	8.910	2.0	ng	27536	1	7.675	272577	20.0
Bicyclo[3.1.0]h...	9.169	6.0	ng	89736	2	10.469	301368	20.0
unknown14.163	14.163	6.6	ng	150945	3	14.339	456213	20.0
Pentane, 2,3,3-...	14.239	5.5	ng	126702	3	14.339	456213	20.0
Silane, trichlo...	15.198	8.3	ng	188418	3	14.339	456213	20.0
Tridecane	16.069	12.7	ng	344963	4	17.104	543905	20.0
Dodecane, 1-iodo-	16.869	6.3	ng	170038	4	17.104	543905	20.0
17-Norkaur-15-e...	18.292	33.6	ng	915167	4	17.104	543905	20.0
2-Phenanthrenol...	20.310	16.3	ng	471598	5	21.204	577551	20.0
Cyclotetracosane	20.916	15.6	ng	451639	5	21.204	577551	20.0
Benzoic acid, 2...	20.986	8.9	ng	255900	5	21.204	577551	20.0
.DELTA.9-THC TFA	21.839	8.7	ng	250158	5	21.204	577551	20.0