

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061522\
 Data File : BP010746.D
 Acq On : 16 Jun 2022 00:37
 Operator : CG/JU
 Sample : N3354-18
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 MW-62

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-BP060722.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BP010746.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.416	389	396	408	rBV	758480	1206574	14.71%	4.281%
2	7.011	657	667	680	rBV	517716	893584	10.89%	3.170%
3	7.828	799	806	815	rBV	223148	366211	4.46%	1.299%
4	8.199	863	869	881	rBV	82973	141467	1.72%	0.502%
5	8.993	997	1004	1020	rVB	1097926	1973127	24.05%	7.000%
6	9.457	1078	1083	1091	rVB	71229	112013	1.37%	0.397%
7	10.628	1270	1282	1287	rBV	325082	568507	6.93%	2.017%
8	13.093	1691	1701	1714	rBV	2925098	4415684	53.83%	15.666%
9	14.192	1882	1888	1896	rBV	78754	122424	1.49%	0.434%
10	14.469	1926	1935	1941	rBV	579709	908086	11.07%	3.222%
11	14.534	1941	1946	1953	rVB	52815	89419	1.09%	0.317%
12	15.639	2126	2134	2138	rBV2	108209	172277	2.10%	0.611%
13	15.963	2182	2189	2198	rBV	3110563	4401466	53.66%	15.616%
14	17.222	2398	2403	2410	rBV	805413	1115644	13.60%	3.958%
15	19.786	2833	2839	2850	rBV	7130091	8203243	100.00%	29.103%
16	21.022	3045	3049	3057	rBV	373328	472280	5.76%	1.676%
17	21.310	3093	3098	3106	rBV	1186285	1508718	18.39%	5.353%
18	23.704	3499	3505	3515	rVB	739433	1515759	18.48%	5.378%

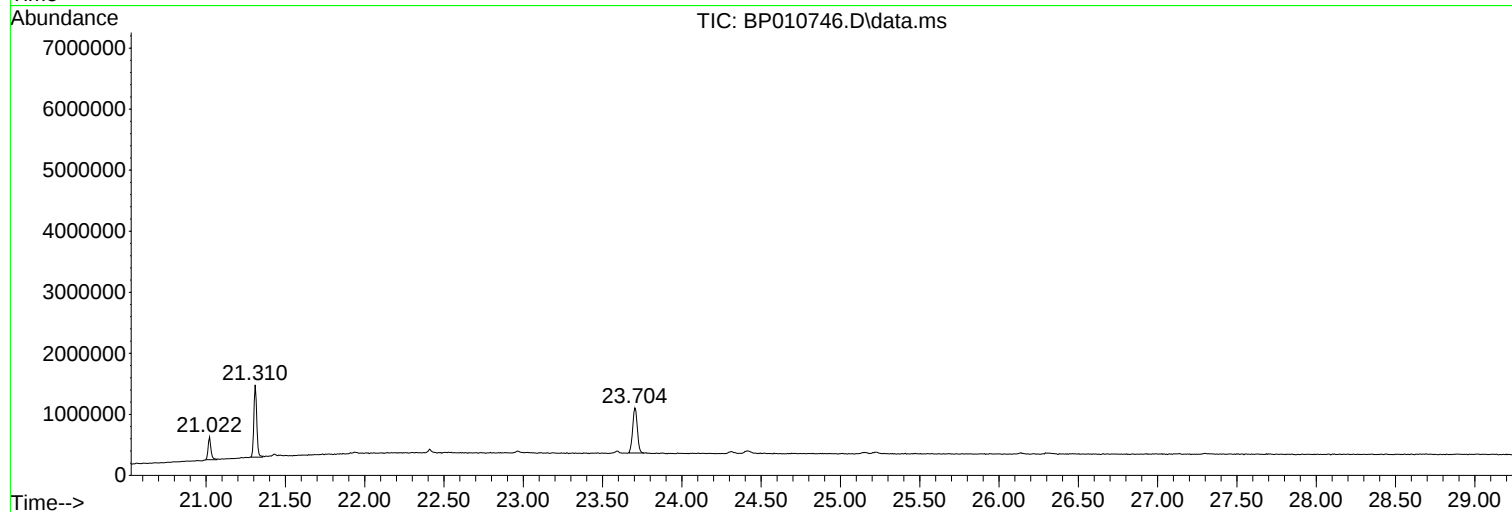
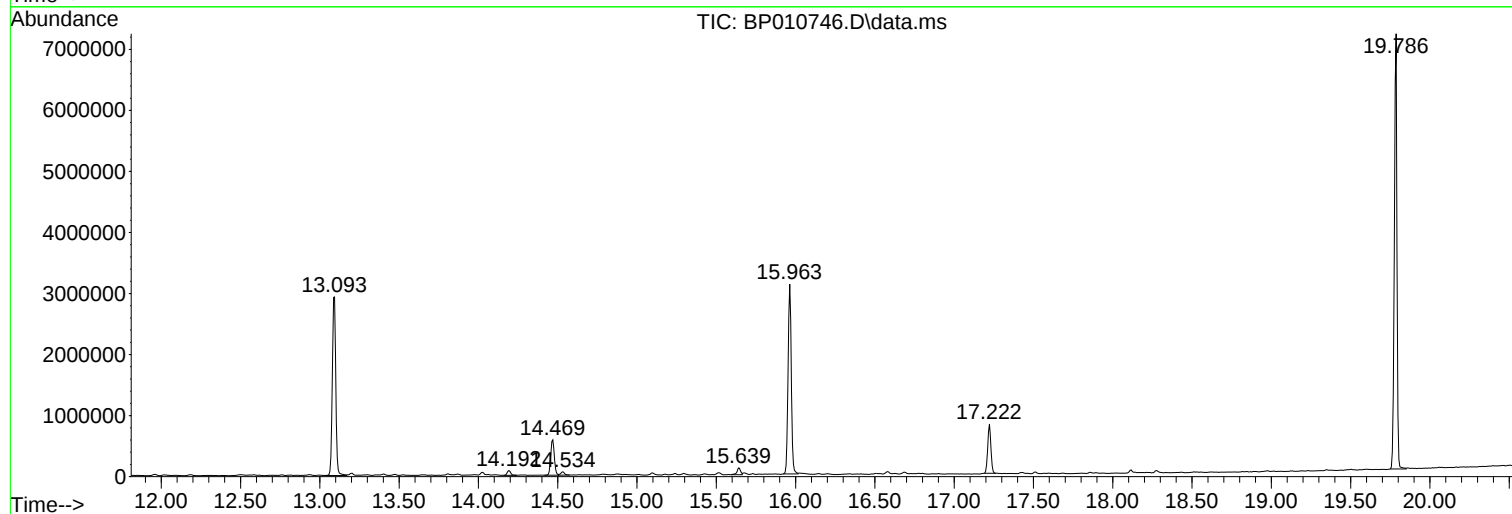
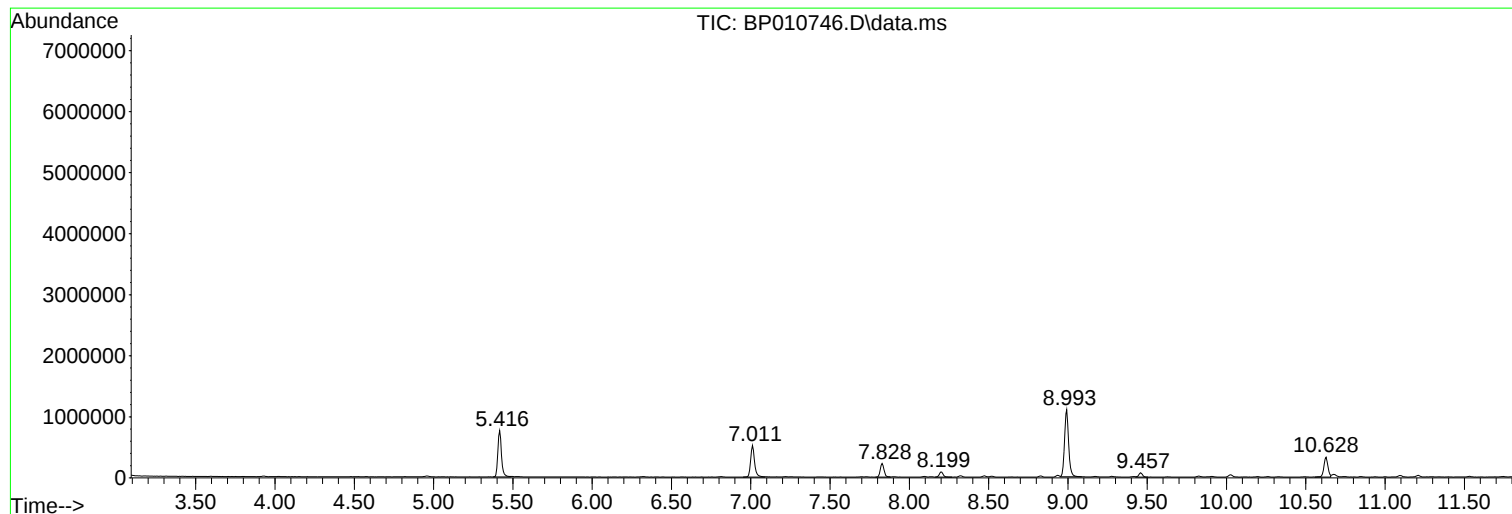
Sum of corrected areas: 28186483

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061522\
Data File : BP010746.D
Acq On : 16 Jun 2022 00:37
Operator : CG/JU
Sample : N3354-18
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
MW-62

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP060722.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061522\
Data File : BP010746.D
Acq On : 16 Jun 2022 00:37
Operator : CG/JU
Sample : N3354-18
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
MW-62

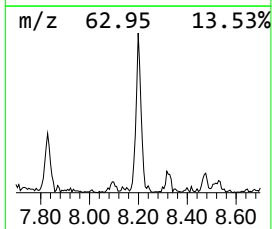
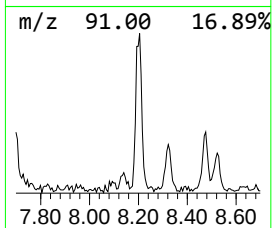
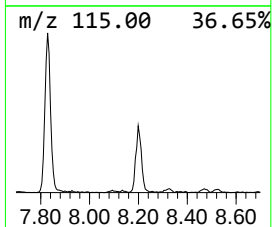
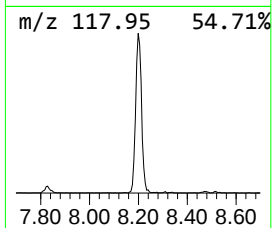
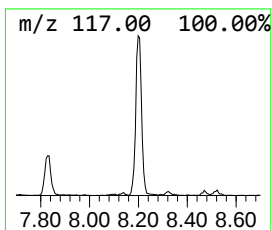
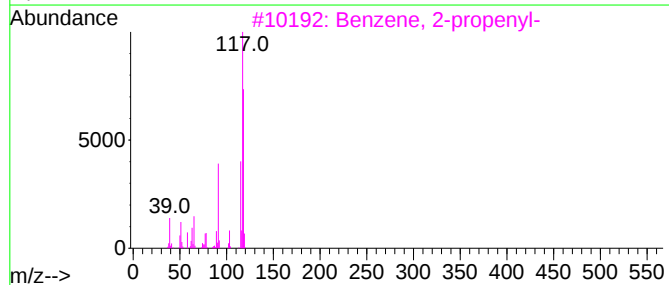
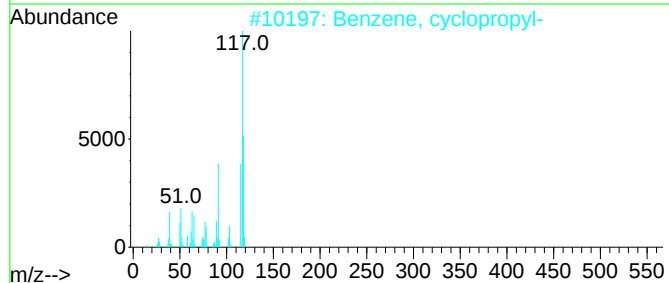
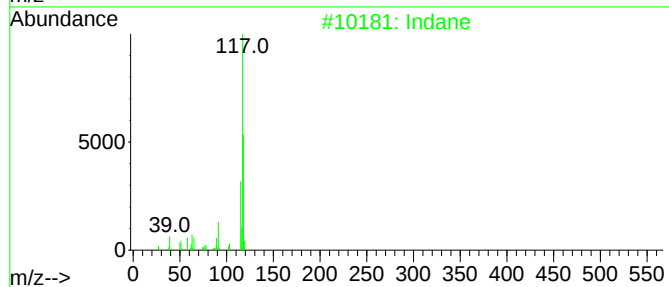
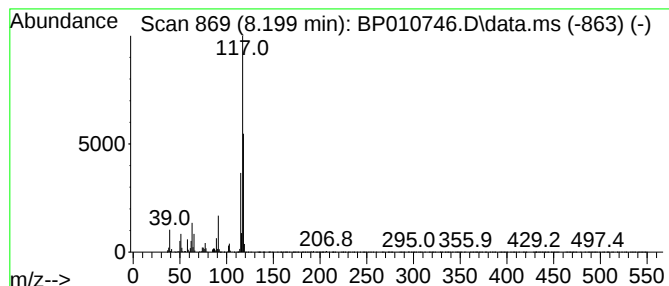
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP060722.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Indane Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.199	7.73 ng	141467	1,4-Dichlorobenzene-d4	7.828

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Indane	118	C9H10	000496-11-7	93
2			Benzene, cyclopropyl-	118	C9H10	000873-49-4	87
3			Benzene, 2-propenyl-	118	C9H10	000300-57-2	72
4			Benzene, 1,1'-(1-ethenyl-1,3-pro...	222	C17H18	061141-97-7	72
5			Tetracyclo[3.3.1.0(2,8).0(4,6)]-...	118	C9H10	1000191-13-7	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061522\
Data File : BP010746.D
Acq On : 16 Jun 2022 00:37
Operator : CG/JU
Sample : N3354-18
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
MW-62

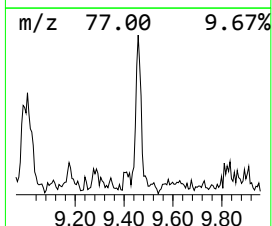
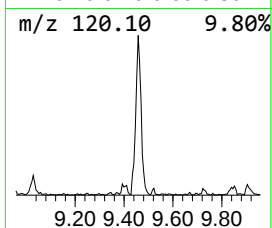
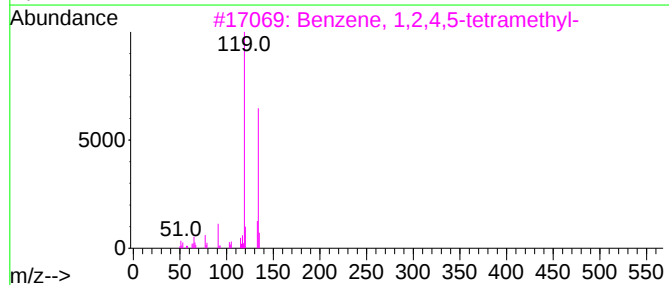
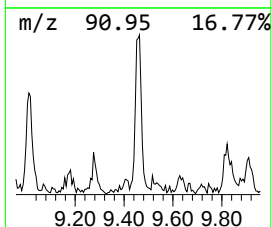
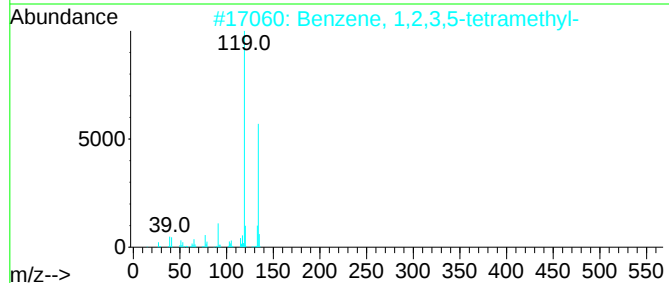
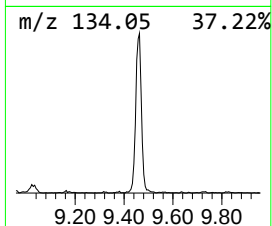
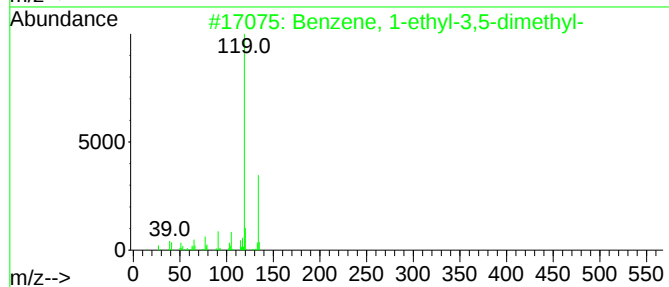
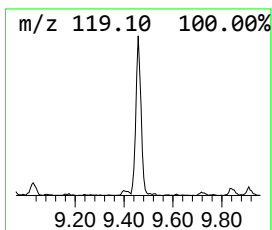
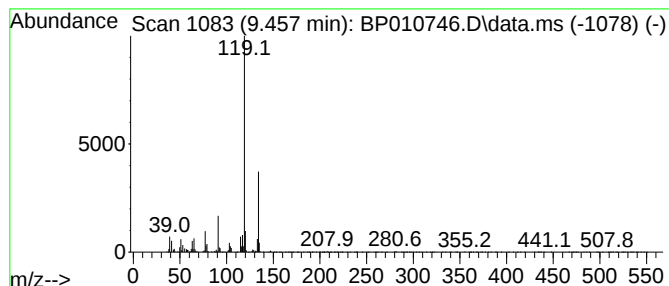
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP060722.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Benzene, 1-ethyl-3,5-dimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.457	3.94 ng	112013	Naphthalene-d8	10.628

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	94
2			Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	94
3			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	94
4			p-Cymene	134	C10H14	000099-87-6	94
5			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061522\
Data File : BP010746.D
Acq On : 16 Jun 2022 00:37
Operator : CG/JU
Sample : N3354-18
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
MW-62

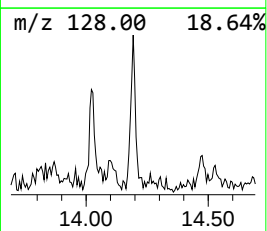
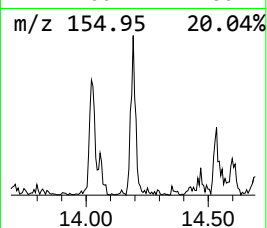
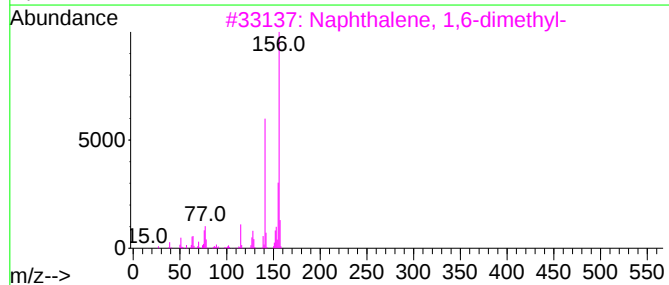
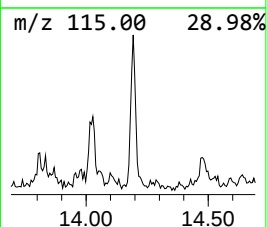
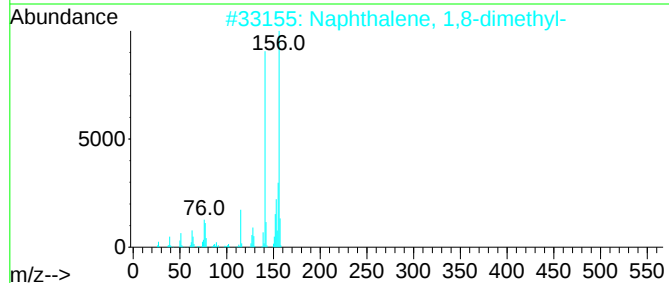
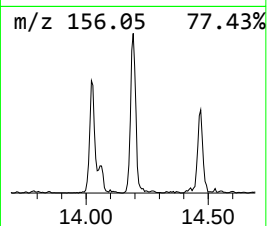
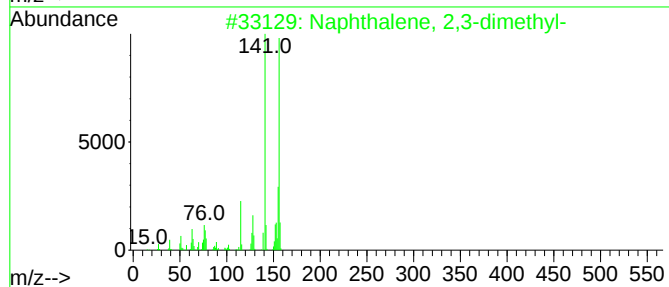
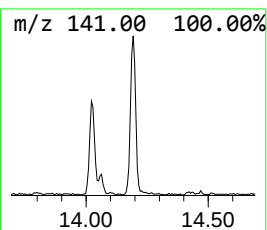
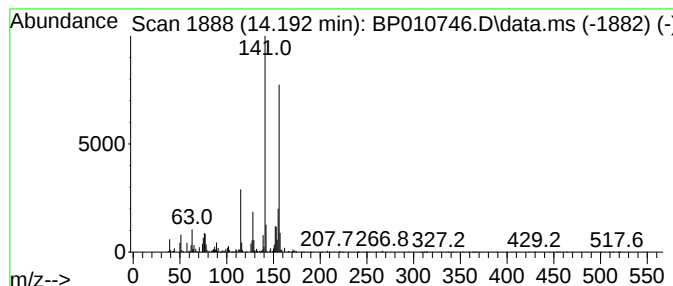
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP060722.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Naphthalene, 2,3-dimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.193	2.70 ng	122424	Acenaphthene-d10	14.469

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	96
2			Naphthalene, 1,8-dimethyl-	156	C12H12	000569-41-5	94
3			Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	94
4			Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	94
5			Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061522\
Data File : BP010746.D
Acq On : 16 Jun 2022 00:37
Operator : CG/JU
Sample : N3354-18
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
MW-62

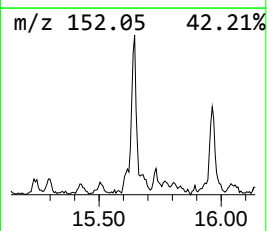
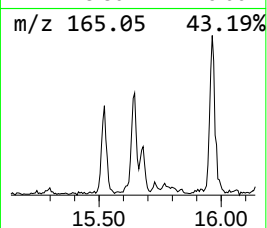
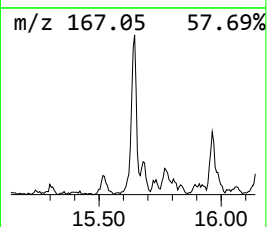
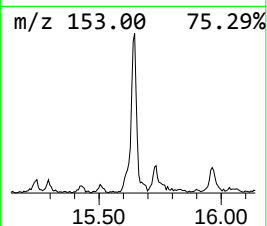
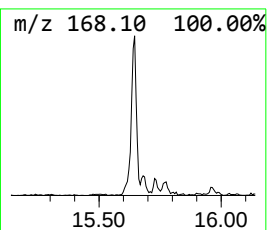
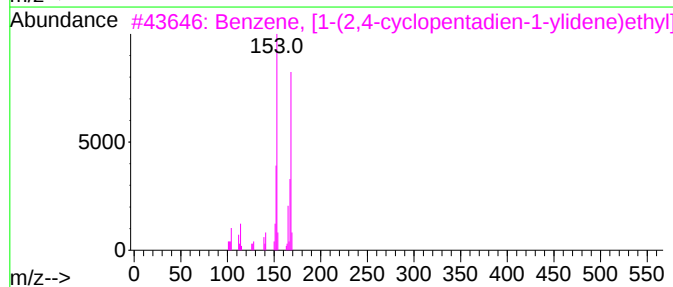
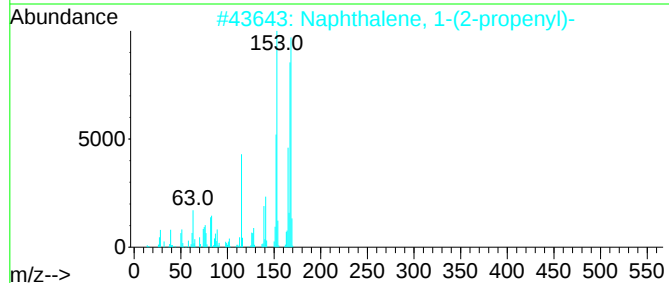
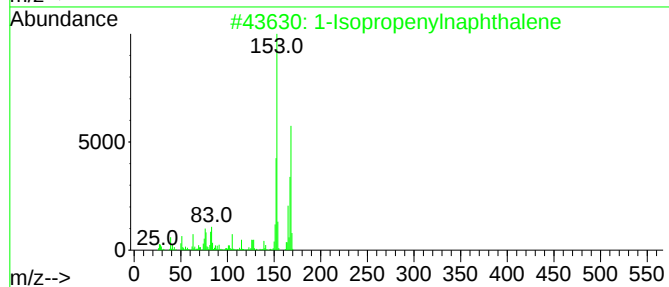
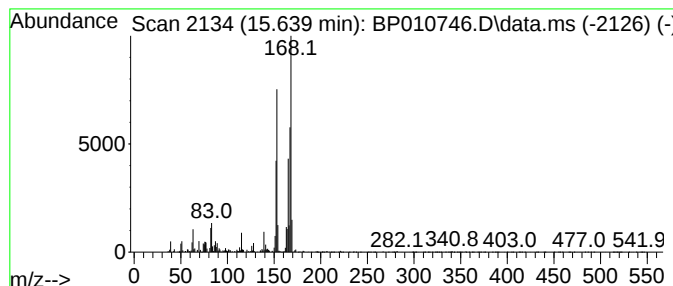
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP060722.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 1-Isopropenylnaphthalene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.640	3.79 ng	172277	Acenaphthene-d10	14.469

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		1-Isopropenylnaphthalene	168	C13H12	001855-47-6	76
2		Naphthalene, 1-(2-propenyl)-	168	C13H12	002489-86-3	62
3		Benzene, [1-(2,4-cyclopentadien-...	168	C13H12	002320-32-3	50
4		1,1'-Biphenyl, 3-methyl-	168	C13H12	000643-93-6	50
5		1,1'-Biphenyl, 2-methyl-	168	C13H12	000643-58-3	46



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061522\
Data File : BP010746.D
Acq On : 16 Jun 2022 00:37
Operator : CG/JU
Sample : N3354-18
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
MW-62

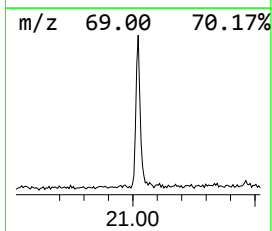
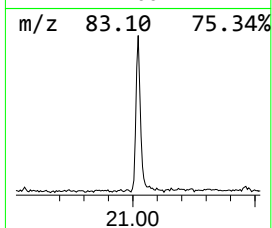
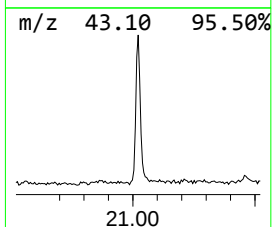
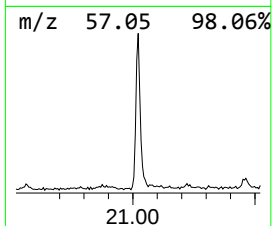
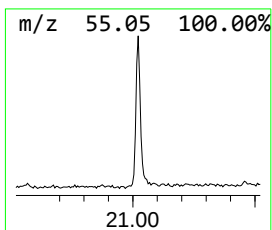
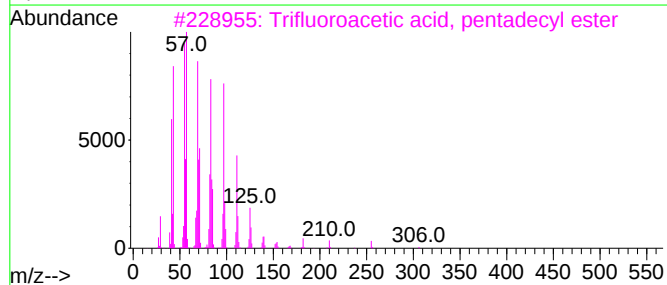
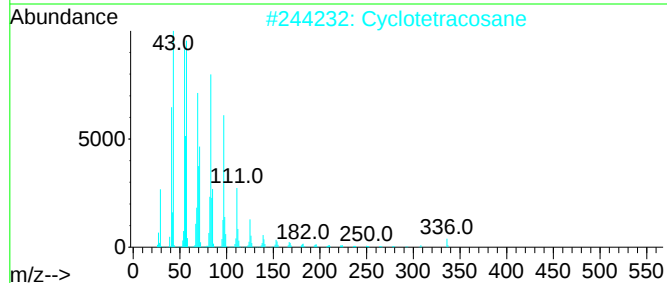
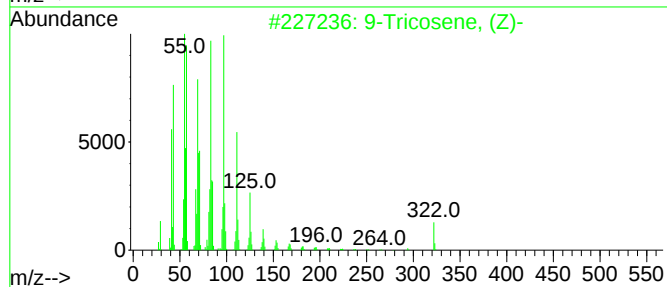
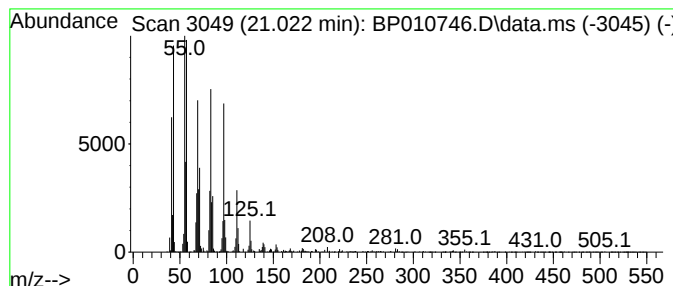
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP060722.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 9-Tricosene, (Z)- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.022	6.26 ng	472280	Chrysene-d12	21.310

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			9-Tricosene, (Z)-	322	C23H46	027519-02-4	96
2			Cyclotetracosane	336	C24H48	000297-03-0	95
3			Trifluoroacetic acid, pentadecyl...	324	C17H31F3O2	959010-23-2	94
4			Trifluoroacetoxy hexadecane	338	C18H33F3O2	006222-03-3	94
5			Pentafluoropropionic acid, penta...	374	C18H31F5O2	959092-08-1	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061522\
Data File : BP010746.D
Acq On : 16 Jun 2022 00:37
Operator : CG/JU
Sample : N3354-18
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
MW-62

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP060722.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Indane	8.199	7.7	ng	141467	1	7.828	366211	20.0
Benzene, 1-ethy...	9.457	3.9	ng	112013	2	10.628	568507	20.0
Naphthalene, 2,...	14.193	2.7	ng	122424	3	14.469	908086	20.0
1-Isopropenylna...	15.640	3.8	ng	172277	3	14.469	908086	20.0
9-Tricosene, (Z)-	21.022	6.3	ng	472280	5	21.310	1508720	20.0