

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG102722\
 Data File : BG055332.D
 Acq On : 27 Oct 2022 20:46
 Operator : CG/JU
 Sample : N5282-01
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SUMP-STONE

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_G\Methods\8270-BG102422.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BG055332.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.315	339	345	354	rBV	77771	129554	3.22%	0.694%
2	5.797	420	427	438	rBV	540731	935847	23.25%	5.015%
3	7.419	694	703	717	rVB	564110	1001365	24.88%	5.366%
4	8.271	835	848	860	rBV	163803	293721	7.30%	1.574%
5	9.446	1040	1048	1061	rBV	359785	653500	16.24%	3.502%
6	11.103	1324	1330	1341	rVB	224953	407947	10.14%	2.186%
7	12.184	1509	1514	1520	rVB4	54263	107327	2.67%	0.575%
8	12.871	1627	1631	1636	rBV6	49452	112871	2.80%	0.605%
9	12.977	1645	1649	1653	rBV	133957	204275	5.08%	1.095%
10	13.095	1666	1669	1677	rBV8	63182	161159	4.00%	0.864%
11	13.406	1718	1722	1729	rVB2	218714	311351	7.74%	1.668%
12	13.512	1735	1740	1745	rVB	893646	1352692	33.61%	7.248%
13	13.676	1764	1768	1776	rBV5	278079	580945	14.43%	3.113%
14	14.311	1872	1876	1879	rBV	717191	1176250	29.22%	6.303%
15	14.346	1879	1882	1886	rVB2	886302	1183593	29.41%	6.342%
16	14.399	1887	1891	1895	rVV3	442760	772191	19.18%	4.138%
17	14.634	1926	1931	1936	rBV5	671828	1491620	37.06%	7.993%
18	14.722	1942	1946	1951	rVB4	1038196	1630639	40.51%	8.738%
19	23.095	3368	3371	3379	rVB	351061	590053	14.66%	3.162%
20	23.635	3458	3463	3479	rVB	674844	1540019	38.26%	8.252%
21	32.401	4938	4955	4985	rVB2	653036	4024990	100.00%	21.568%

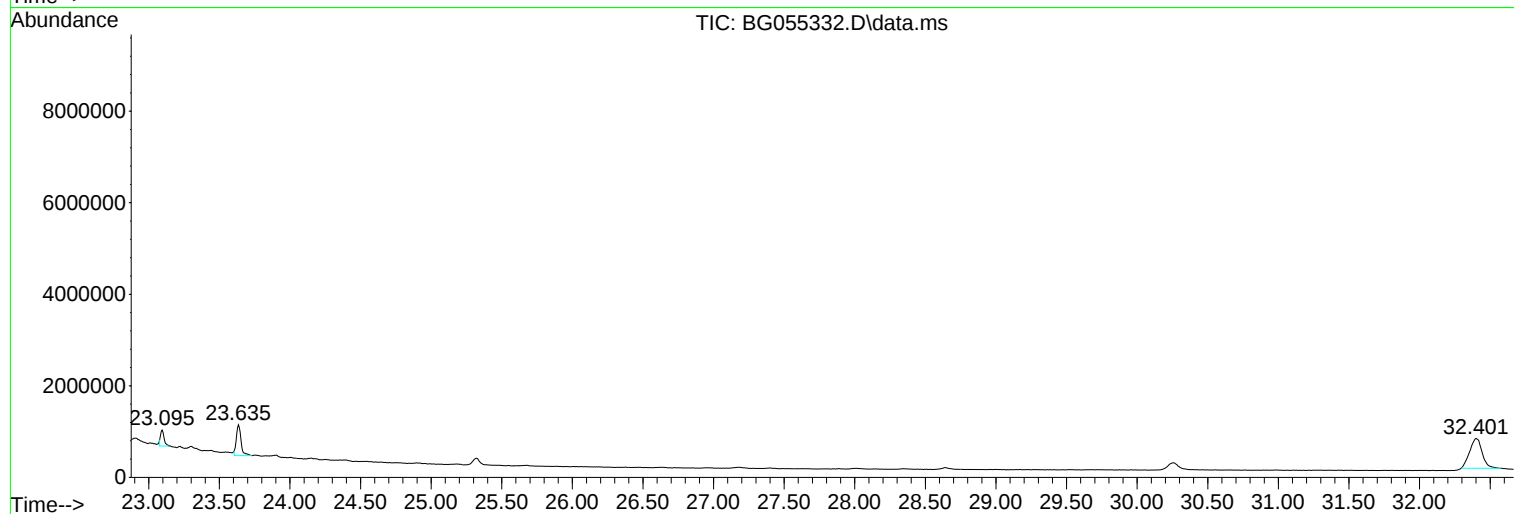
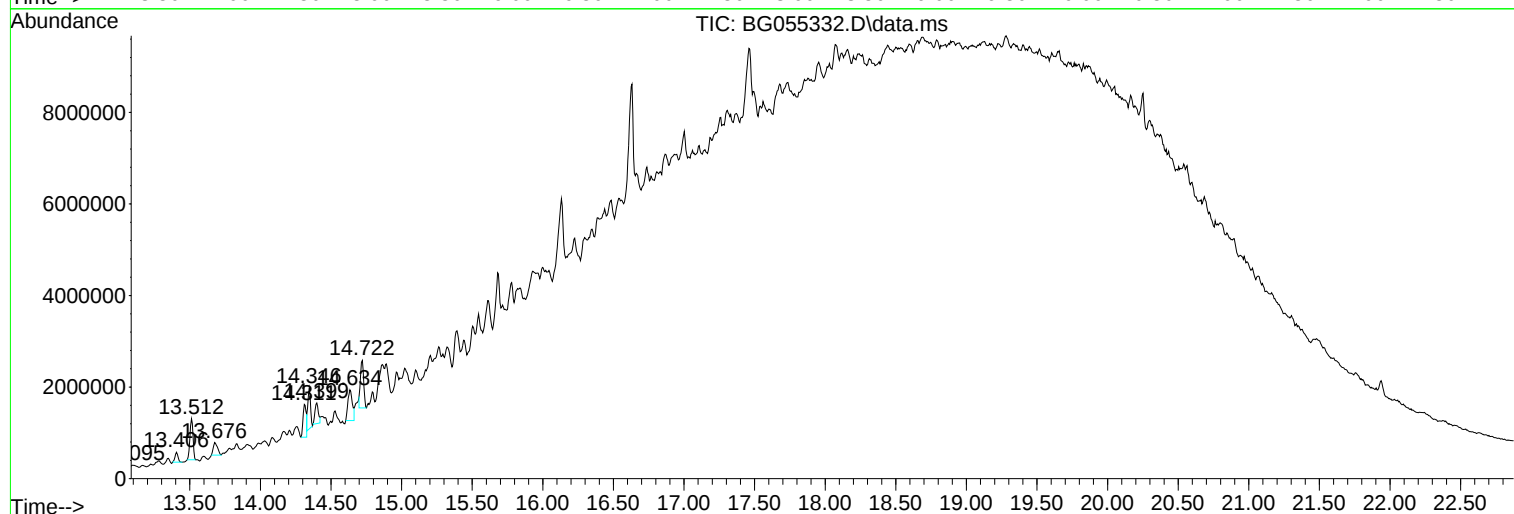
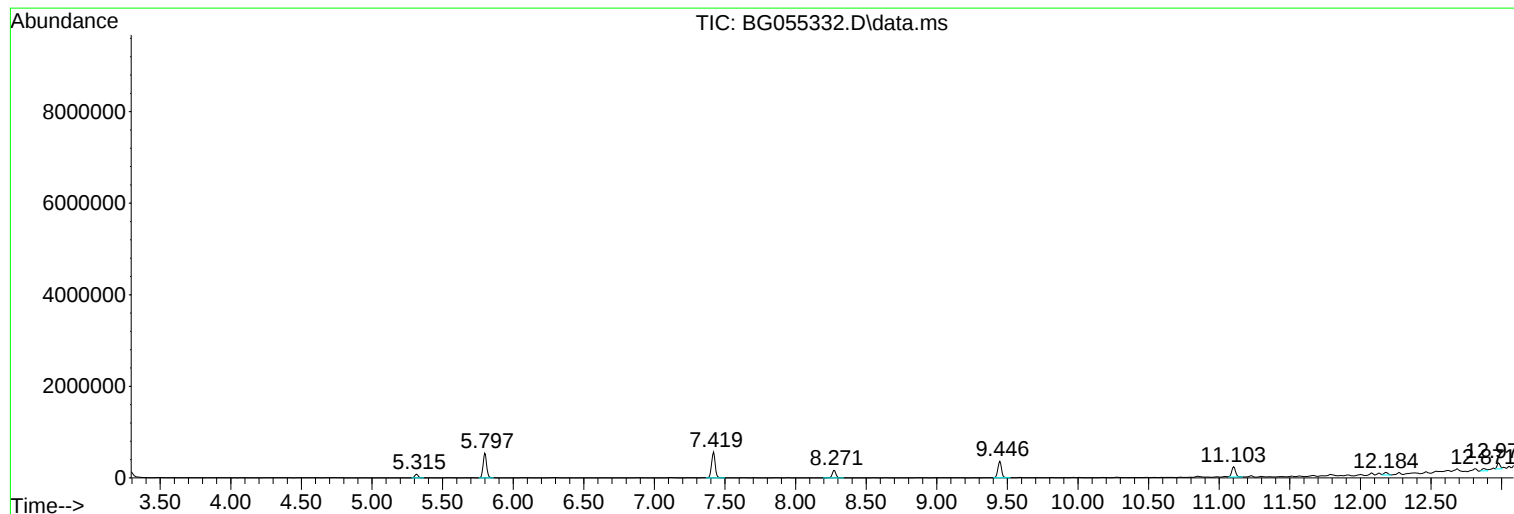
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG102422.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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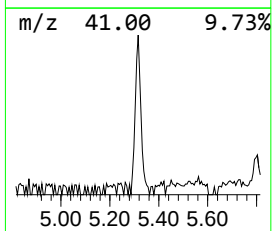
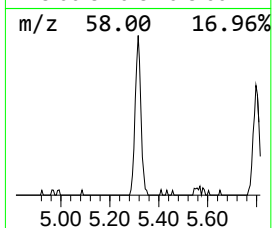
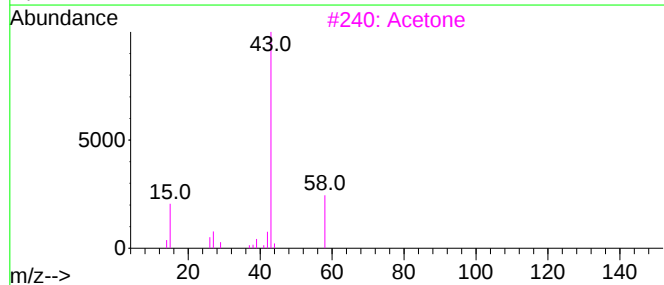
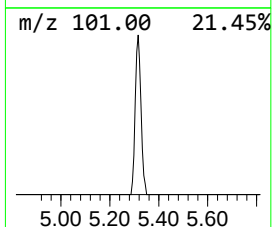
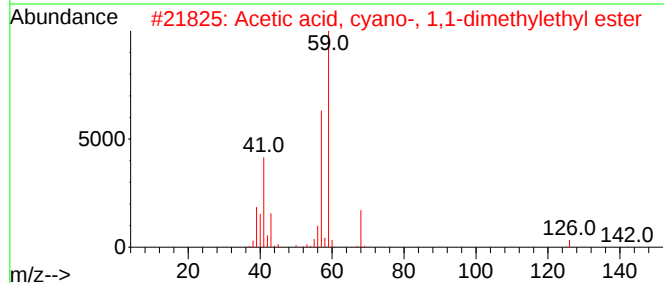
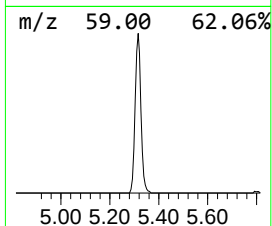
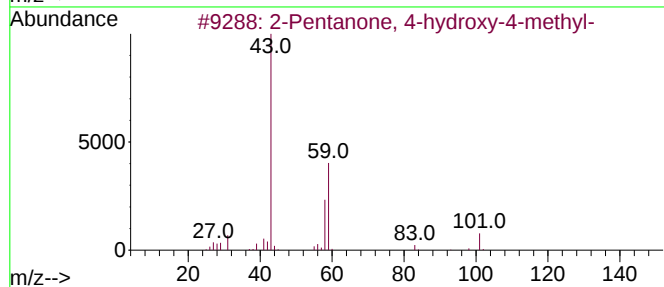
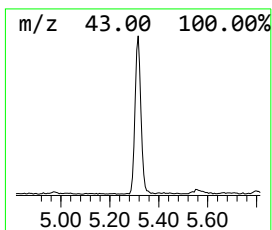
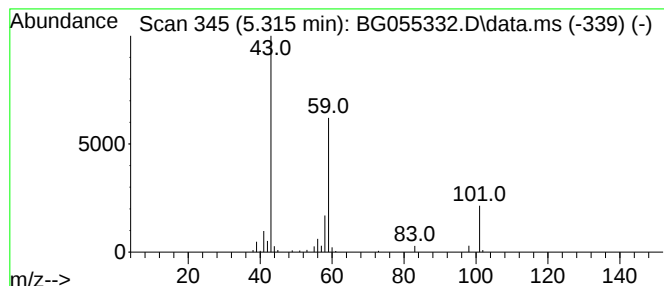
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.315	8.82 ng	129554	1,4-Dichlorobenzene-d4	8.271

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56		
2	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	25		
3	Acetone	58	C3H6O	000067-64-1	9		
4	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9		
5	Propane, 2-ethoxy-2-methyl-	102	C6H14O	000637-92-3	9		



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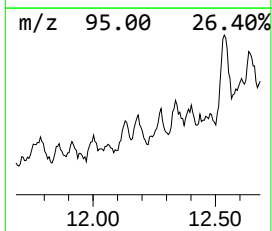
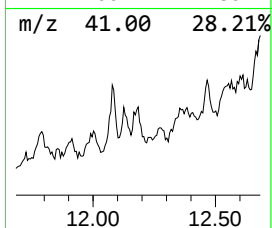
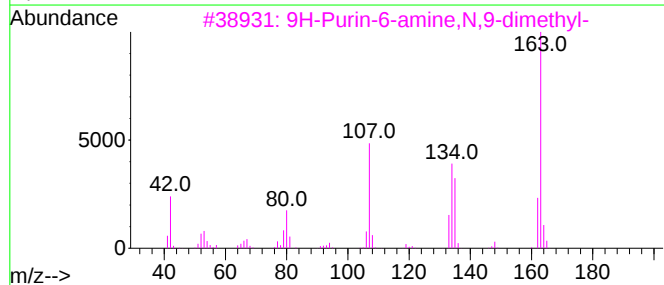
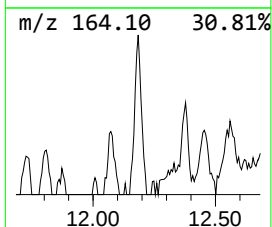
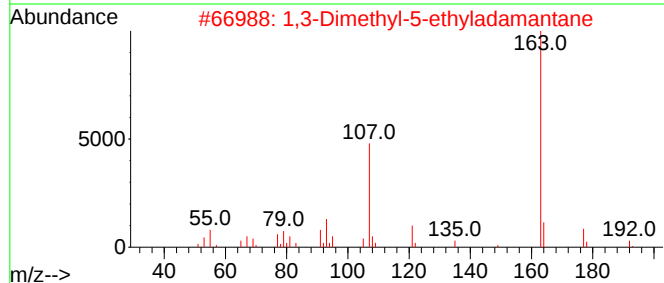
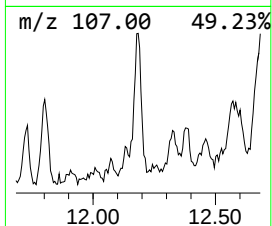
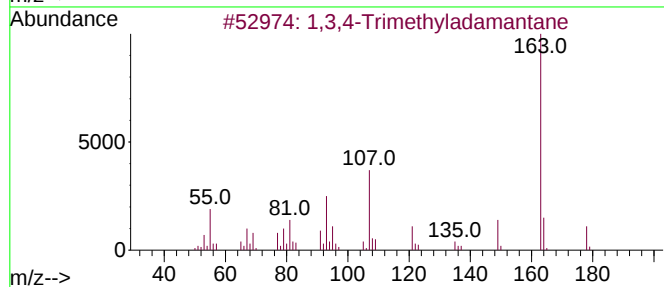
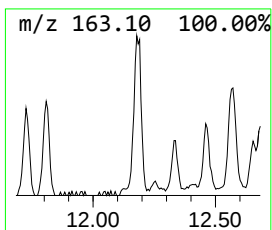
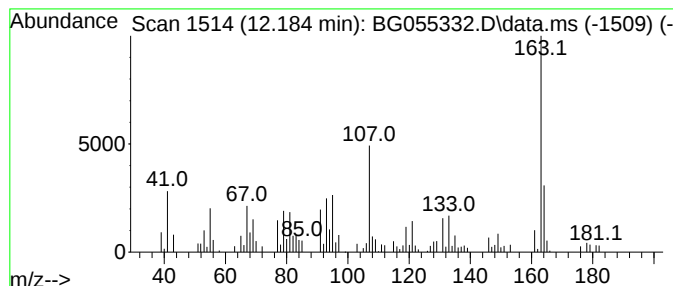
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 1,3,4-Trimethyladamantane Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.184	5.26 ng	107327	Naphthalene-d8	11.103

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,3,4-Trimethyladamantane	178	C13H22	1000214-98-3	53		
2	1,3-Dimethyl-5-ethyladamantane	192	C14H24	001687-35-0	50		
3	9H-Purin-6-amine,N,9-dimethyl-	163	C7H9N5	002009-52-1	49		
4	1,2-Dimethyl-5-nitroadamantane	209	C12H19NO2	006588-68-7	47		
5	1,3,6-Trimethyladamantane	178	C13H22	024139-37-5	42		



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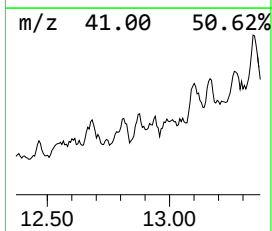
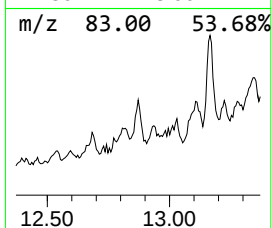
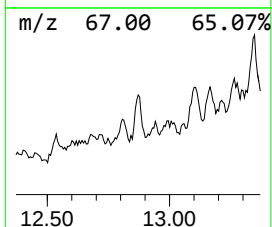
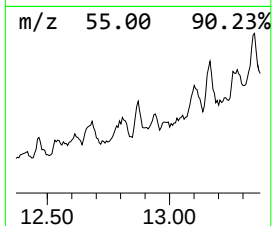
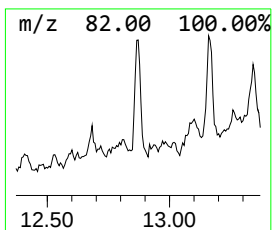
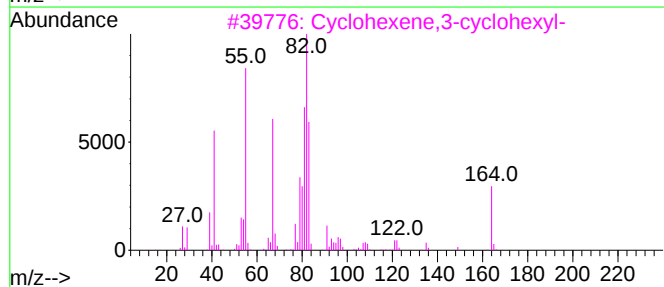
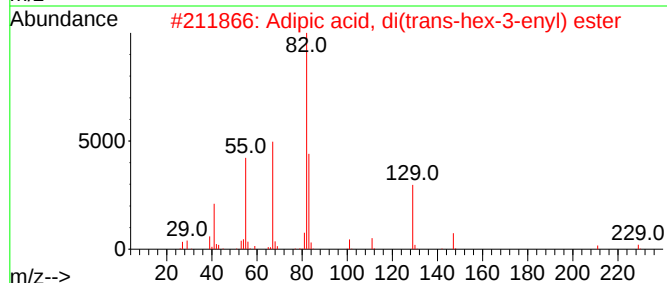
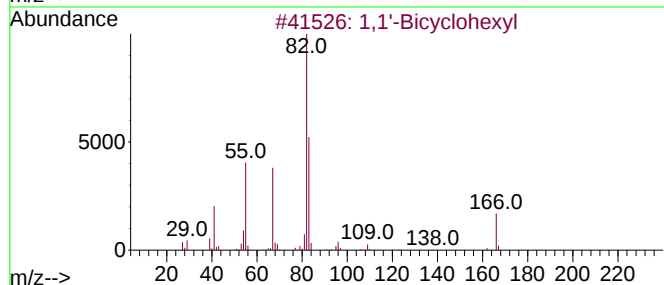
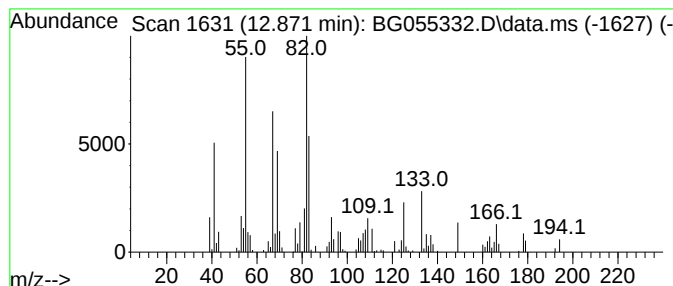
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 1,1'-Bicyclohexyl Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.871	5.53 ng	112871	Naphthalene-d8	11.103

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,1'-Bicyclohexyl	166	C12H22	000092-51-3	53
2			Adipic acid, di(trans-hex-3-enyl...)	310	C18H30O4	1000354-02-3	50
3			Cyclohexene,3-cyclohexyl-	164	C12H20	001808-09-9	50
4			2,4-Hexadiene	82	C6H10	000592-46-1	43
5			Succinic acid, cyclohexylmethyl ...	296	C17H28O4	1000391-08-5	43



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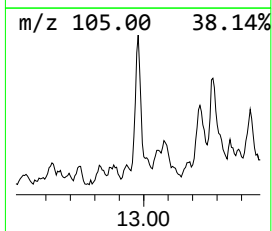
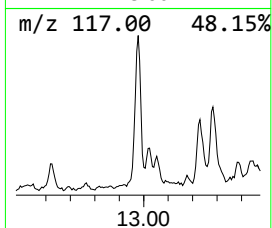
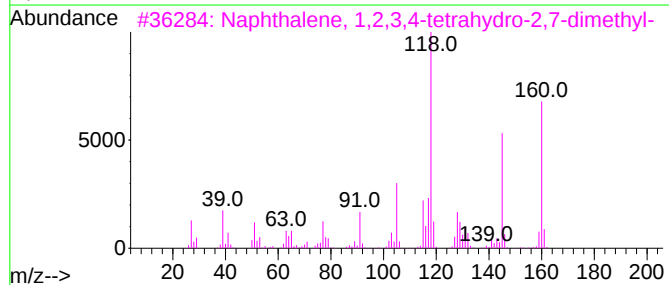
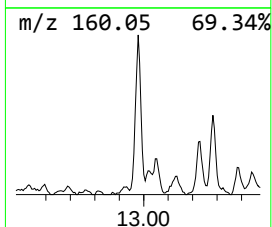
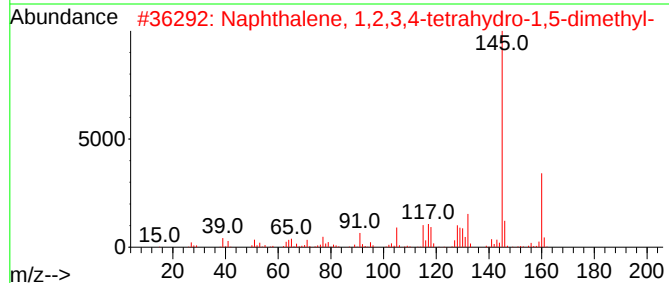
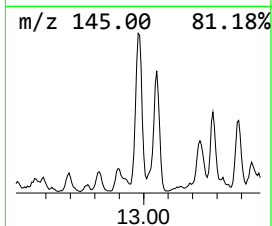
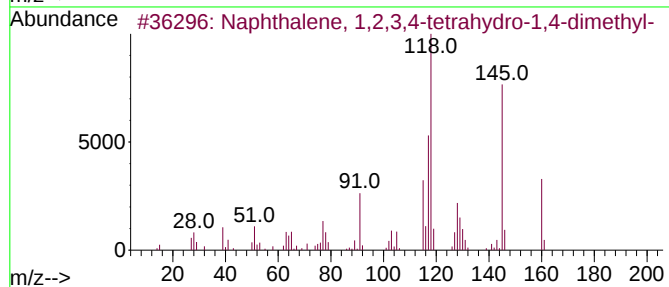
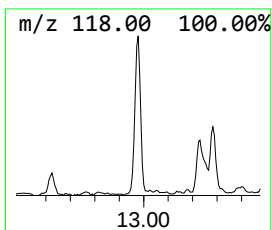
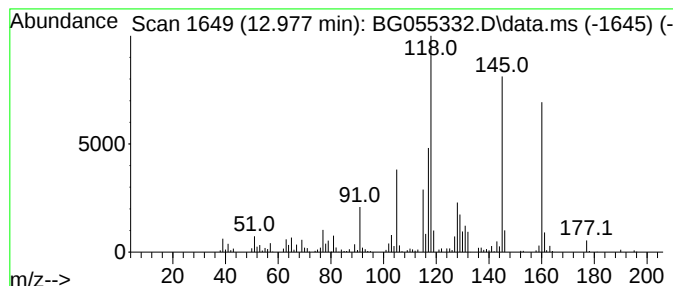
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Peak Number 4 Naphthalene, 1,2,3,4-tetrahydro-... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.977	10.01 ng	204275	Naphthalene-d8	11.103

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	004175-54-6	96
2			Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	021564-91-0	81
3			Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	013065-07-1	74
4			Benzene, 1,3,5-trimethyl-2-(1-me...	160	C12H16	014679-13-1	70
5			Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	007524-63-2	64



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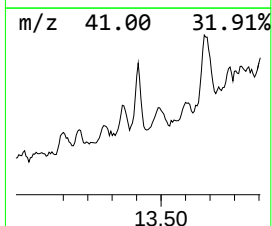
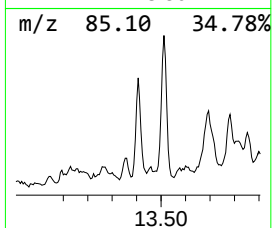
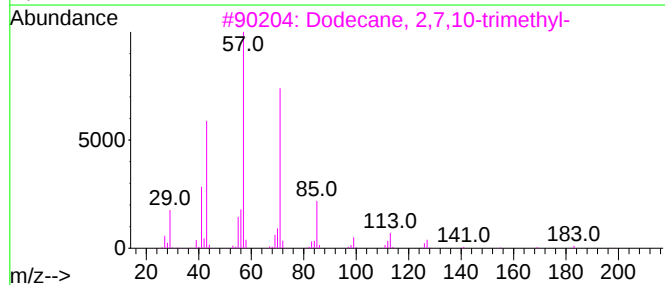
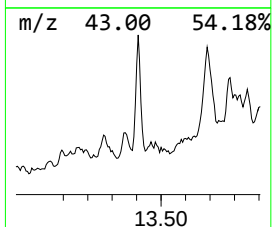
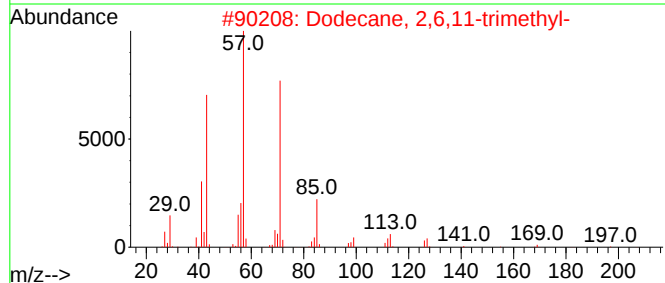
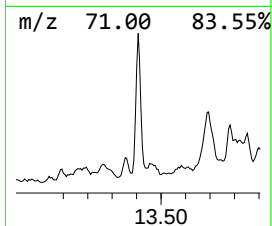
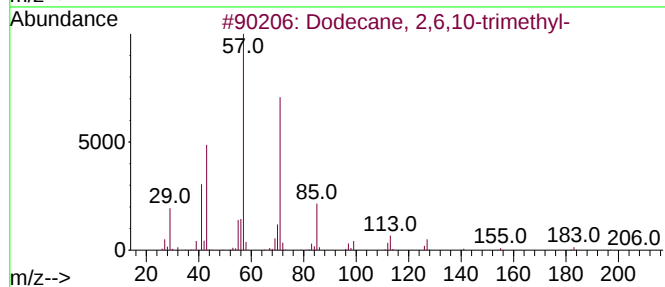
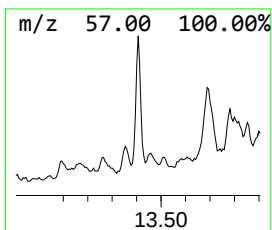
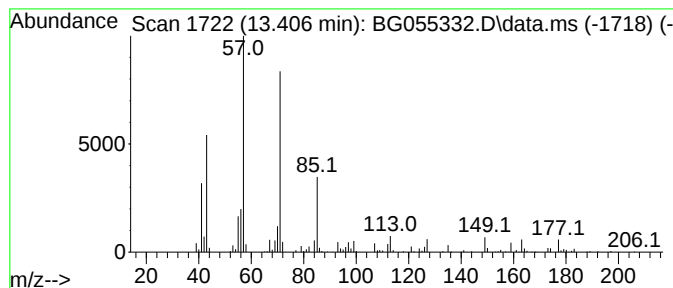
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Dodecane, 2,6,10-trimethyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.406	3.82 ng	311351	Acenaphthene-d10	14.893

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dodecane, 2,6,10-trimethyl-	212	C15H32	003891-98-3	80
2			Dodecane, 2,6,11-trimethyl-	212	C15H32	031295-56-4	72
3			Dodecane, 2,7,10-trimethyl-	212	C15H32	074645-98-0	72
4			Tridecane, 5-propyl-	226	C16H34	055045-11-9	64
5			Sulfurous acid, 2-ethylhexyl non...	320	C17H36O3S	1010309-19-2	53



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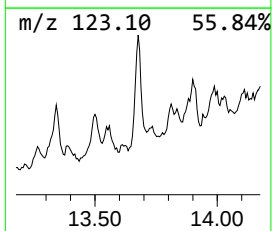
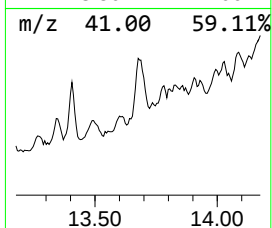
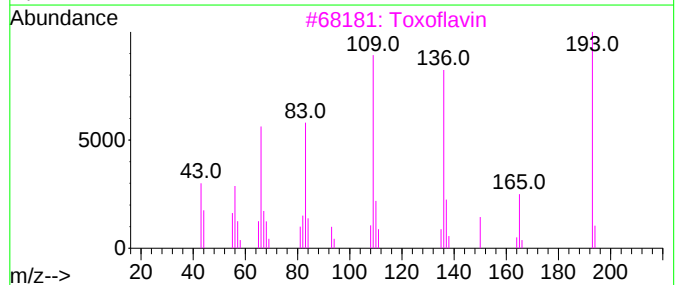
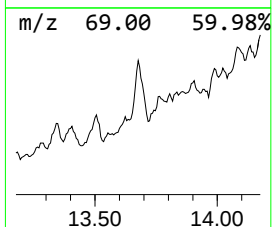
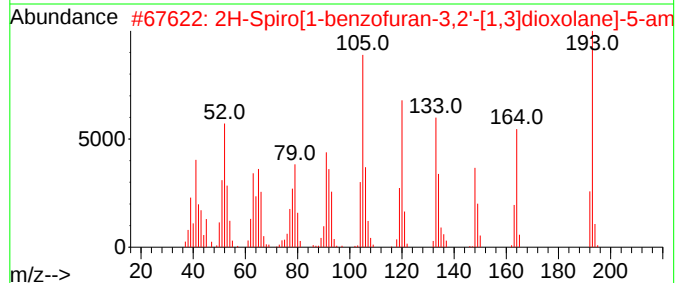
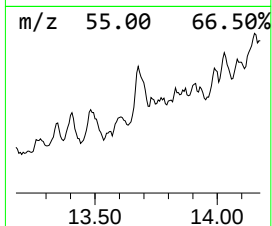
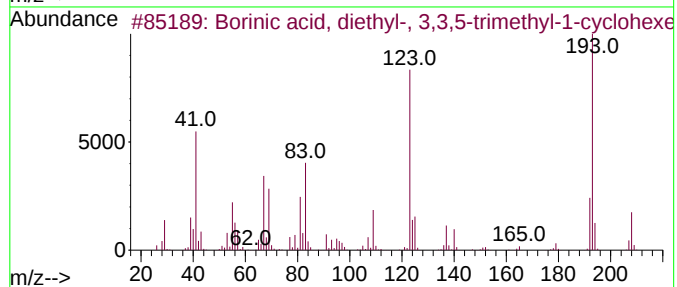
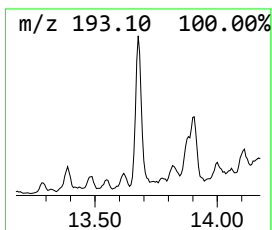
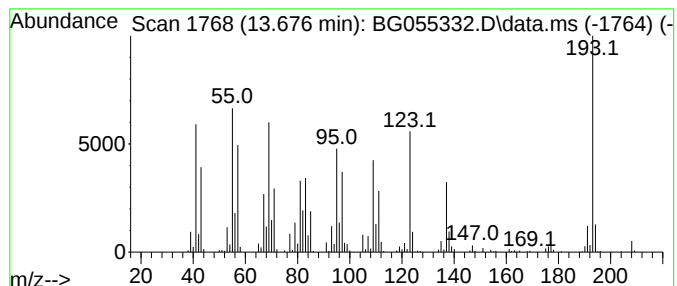
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 unknown13.676 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.676	7.13 ng	580945	Acenaphthene-d10	14.893

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Borinic acid, diethyl-, 3,3,5-tr...	208	C13H25BO	057387-76-5	38
2			2H-Spiro[1-benzofuran-3,2'-[1,3]...	193	C10H11NO3	1000387-33-4	35
3			Toxoflavin	193	C7H7N5O2	000084-82-2	30
4			benzoxazole, 5,6-dimethoxy-2-met...	193	C10H11NO3	1000395-98-9	27
5			9-Vinylcarbazole	193	C14H11N	001484-13-5	25



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG102722\
Data File : BG055332.D
Acq On : 27 Oct 2022 20:46
Operator : CG/JU
Sample : N5282-01
Misc :
ALS Vial : 16 Sample Multiplier: 1

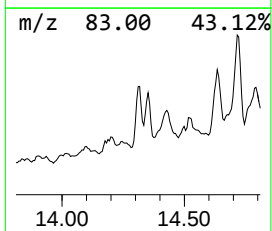
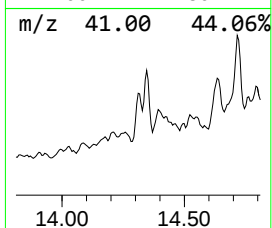
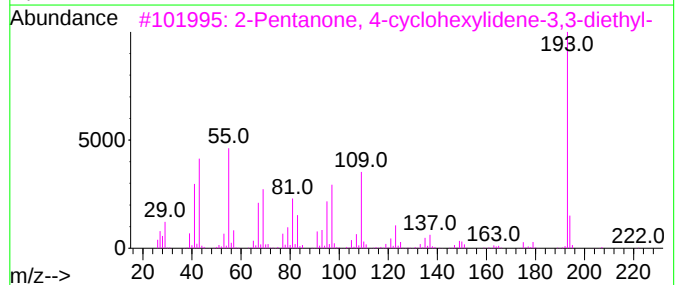
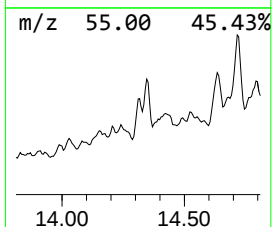
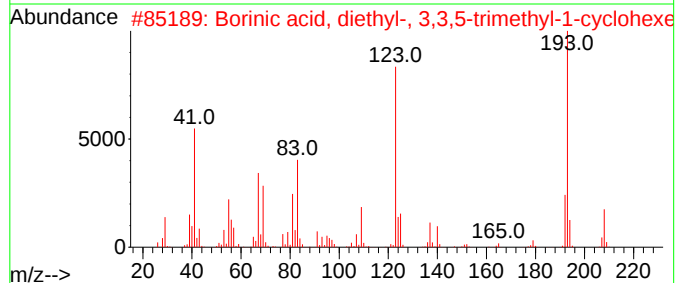
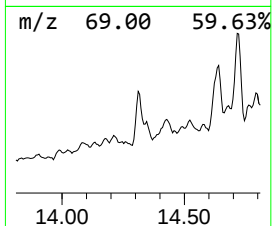
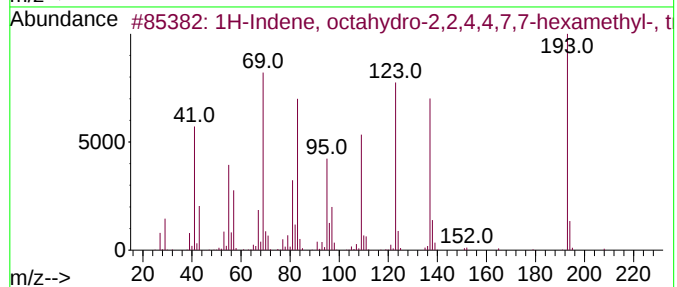
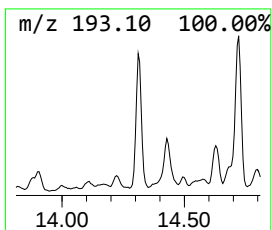
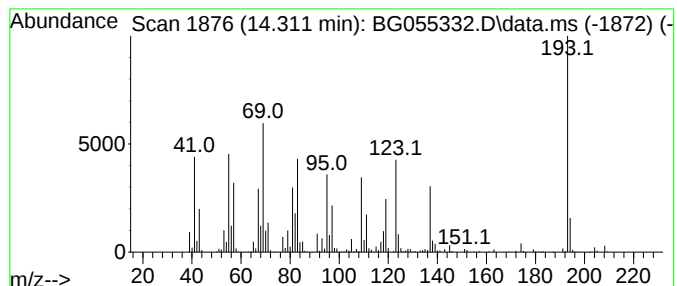
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BNA_G
ClientSampleId :
SUMP-STONE

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG102422.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 1H-Indene, octahydro-2,2,4,... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.311	14.43 ng	1176250	Acenaphthene-d10	14.893	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, octahydro-2,2,4,4,7,7...	208	C15H28	054832-83-6	50
2	Borinic acid, diethyl-, 3,3,5-tr...	208	C13H25BO	057387-76-5	47
3	2-Pentanone, 4-cyclohexylidene-3...	222	C15H26O	313253-65-5	43
4	1-(p-Fluorophenyl)-4-piperidone	193	C11H12FNO	1000238-56-7	43
5	S-Phenyl 5-(phenylthio)pentaneth...	302	C17H18OS2	1000234-40-7	37



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG102722\
Data File : BG055332.D
Acq On : 27 Oct 2022 20:46
Operator : CG/JU
Sample : N5282-01
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SUMP-STONE

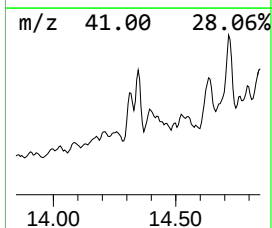
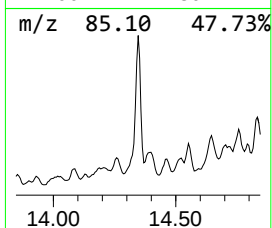
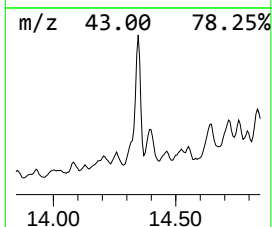
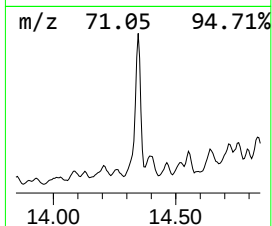
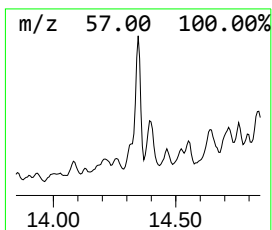
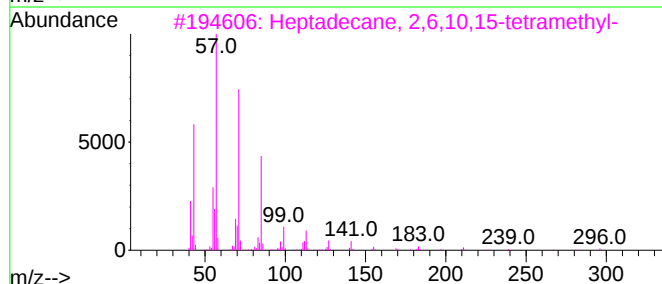
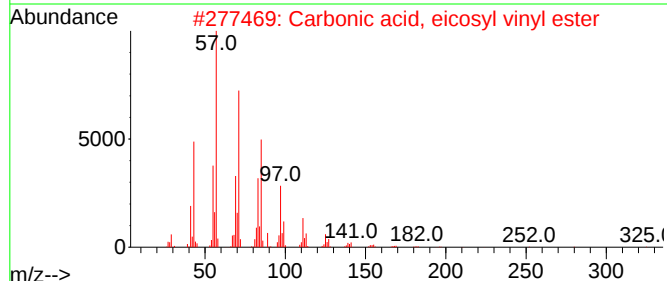
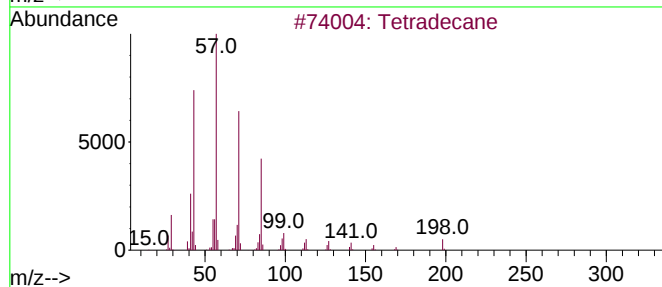
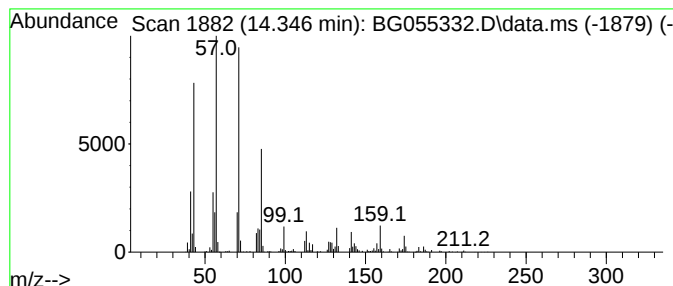
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Tetradecane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.346	14.52 ng	1183590	Acenaphthene-d10	14.893

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetradecane	198	C14H30	000629-59-4	80
2			Carbonic acid, eicosyl vinyl ester	368	C23H44O3	1000382-54-3	80
3			Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	72
4			Decane, 5-propyl-	184	C13H28	017312-62-8	72
5			Nonadecane	268	C19H40	000629-92-5	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG102722\
Data File : BG055332.D
Acq On : 27 Oct 2022 20:46
Operator : CG/JU
Sample : N5282-01
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SUMP-STONE

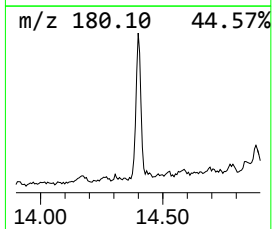
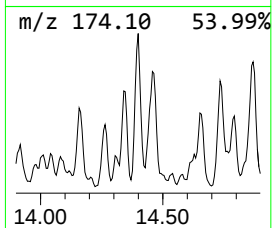
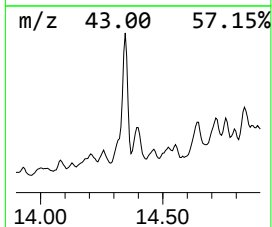
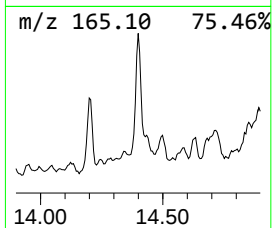
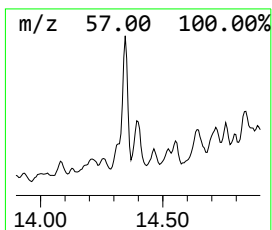
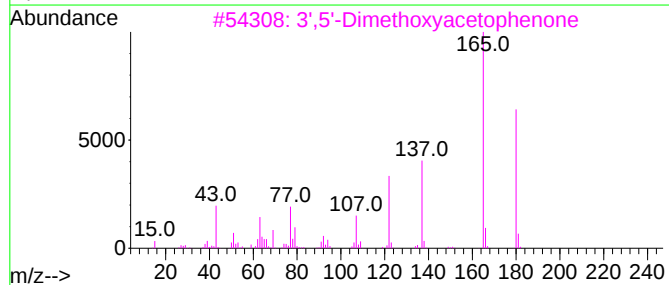
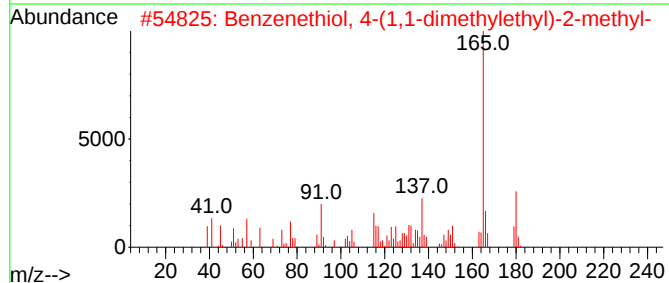
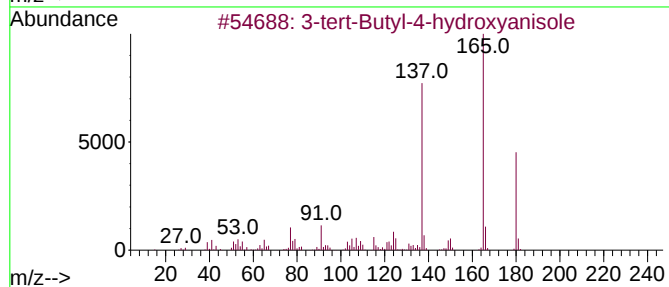
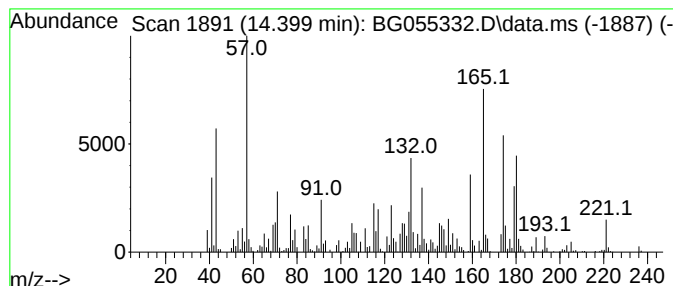
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 3-tert-Butyl-4-hydroxyanisole Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.399	9.47 ng	772191	Acenaphthene-d10	14.893

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		3-tert-Butyl-4-hydroxyanisole	180	C11H16O2	000121-00-6	86
2		Benzenethiol, 4-(1,1-dimethyleth...	180	C11H16S	015570-10-2	49
3		3',5'-Dimethoxyacetophenone	180	C10H12O3	039151-19-4	46
4		o-Cresol, TMS derivative	180	C10H16OSi	001009-02-5	42
5		4a,9a-Methano-9H-fluorene	180	C14H12	019540-84-2	41



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG102722\
Data File : BG055332.D
Acq On : 27 Oct 2022 20:46
Operator : CG/JU
Sample : N5282-01
Misc :
ALS Vial : 16 Sample Multiplier: 1

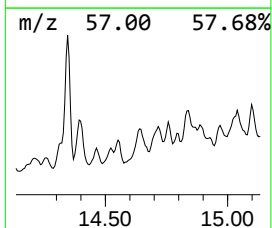
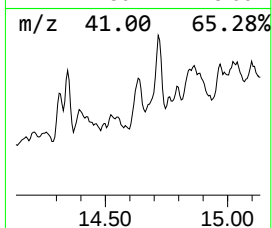
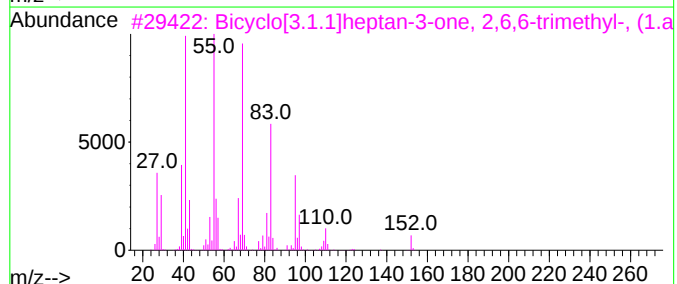
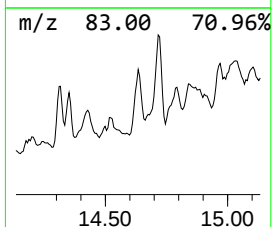
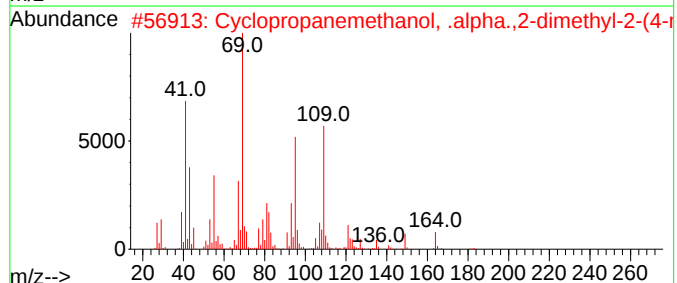
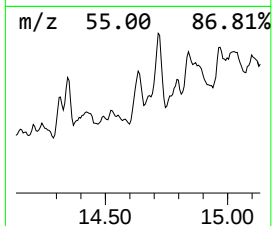
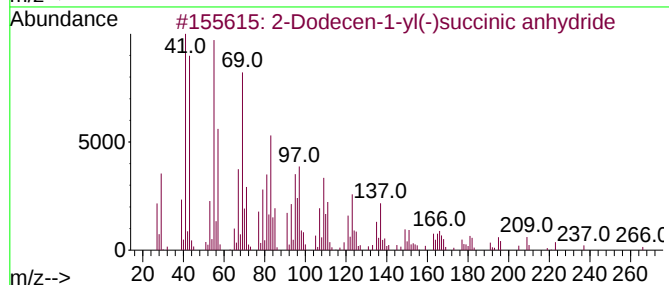
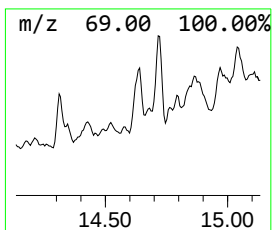
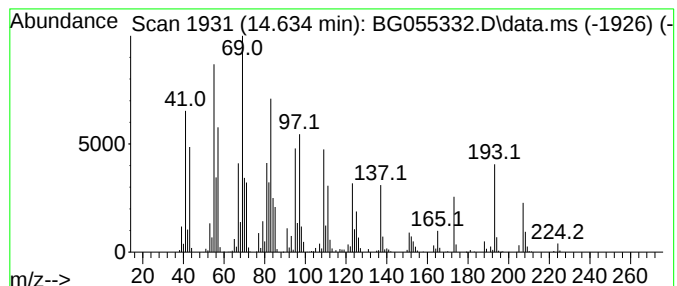
Instrument :
BNA_G
ClientSampleId :
SUMP-STONE

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG102422.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 2-Dodecen-1-yl(-)succinic a... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.634	18.29 ng	1491620	Acenaphthene-d10	14.893	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Dodecen-1-yl(-)succinic anhydride	266	C16H26O3	019780-11-1	58
2	Cyclopropanemethanol, .alpha.,2-...	182	C12H22O	121959-70-4	38
3	Bicyclo[3.1.1]heptan-3-one, 2,6,...	152	C10H16O	015358-88-0	27
4	1,1'-Bicyclohexyl, 2-propyl-, cis-	208	C15H28	054934-88-2	25
5	1,7-Dimethyl-4-(1-methylethyl)cy...	210	C15H30	000645-10-3	22



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG102722\
Data File : BG055332.D
Acq On : 27 Oct 2022 20:46
Operator : CG/JU
Sample : N5282-01
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SUMP-STONE

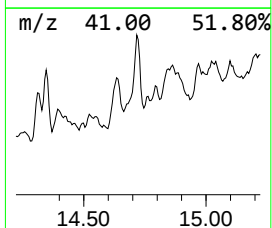
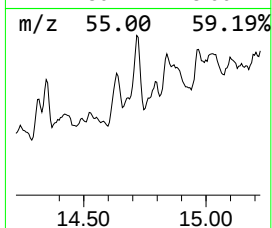
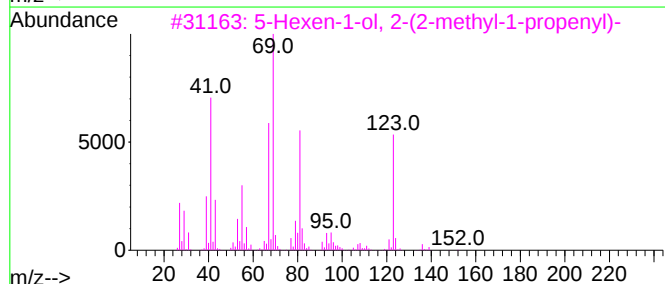
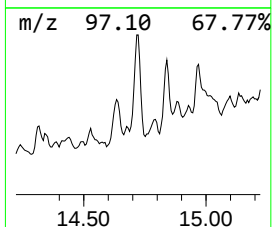
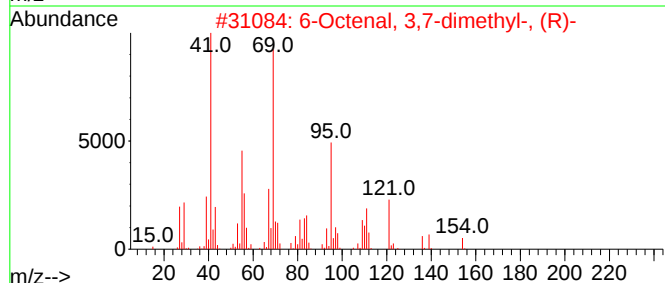
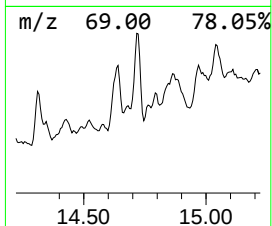
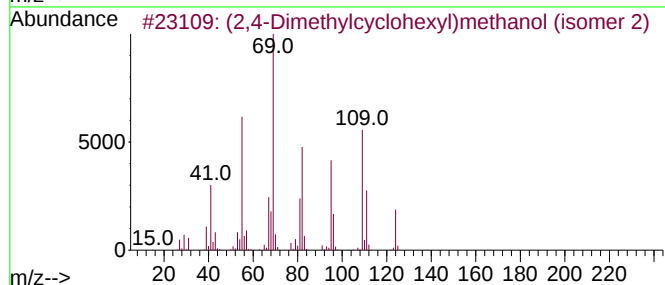
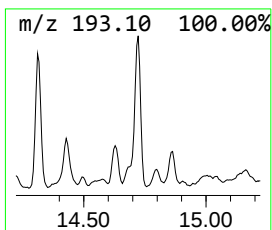
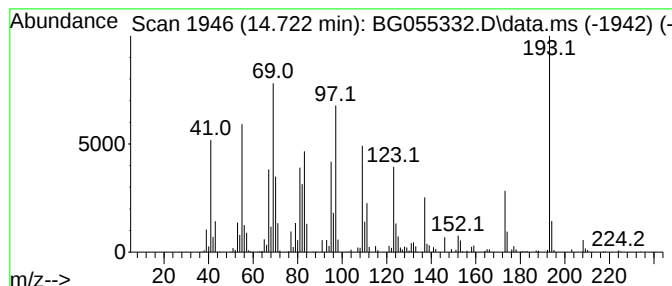
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 unknown14.722 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.722	20.00 ng	1630640	Acenaphthene-d10	14.893

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			(2,4-Dimethylcyclohexyl)methanol...	142	C9H18O	1000499-02-7	27
2			6-Octenal, 3,7-dimethyl-, (R)-	154	C10H18O	002385-77-5	22
3			5-Hexen-1-ol, 2-(2-methyl-1-prop...	154	C10H18O	018675-19-9	14
4			1-Decene, 10-bromo-	218	C10H19Br	062871-09-4	14
5			Silane, chlorodiethylhexyloxy-	222	C10H23ClOSi	1000363-74-4	14



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG102722\
Data File : BG055332.D
Acq On : 27 Oct 2022 20:46
Operator : CG/JU
Sample : N5282-01
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SUMP-STONE

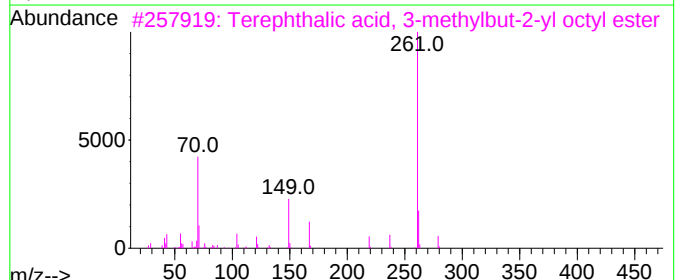
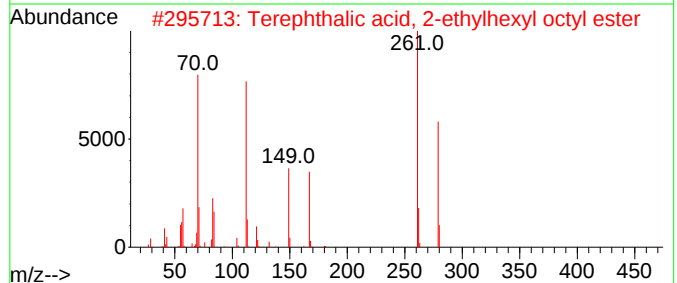
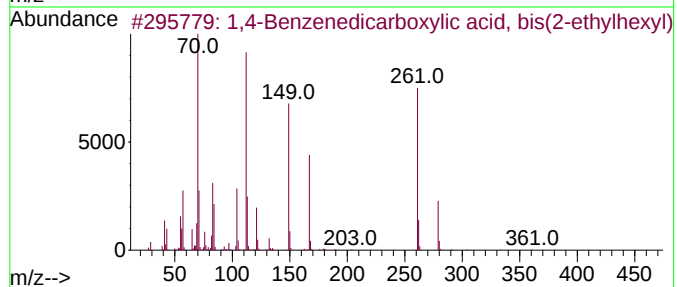
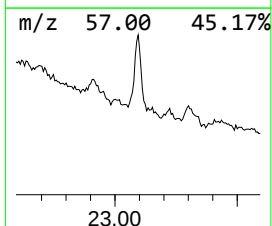
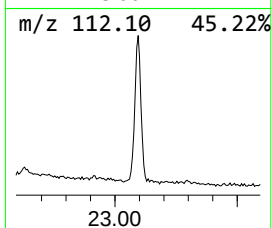
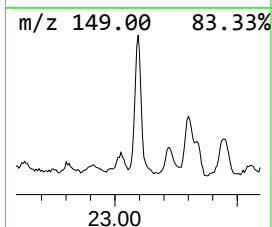
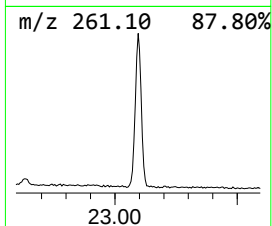
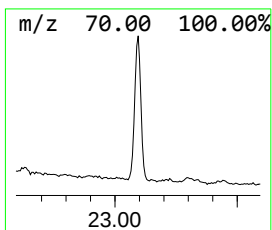
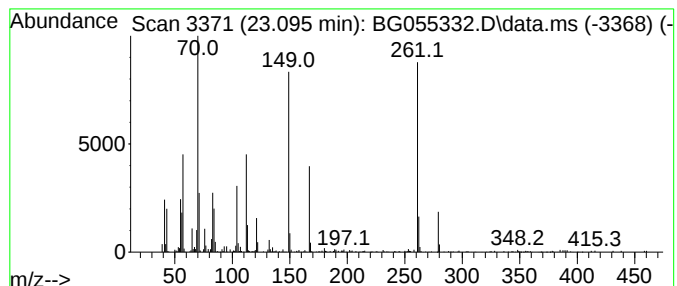
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG102422.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 1,4-Benzenedicarboxylic aci... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.095	20.00 ng	590053	Chrysene-d12	21.937

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,4-Benzenedicarboxylic acid, bi...	390	C24H38O4	006422-86-2	81
2			Terephthalic acid, 2-ethylhexyl ...	390	C24H38O4	1000324-00-5	62
3			Terephthalic acid, 3-methylbut-2...	348	C21H32O4	1000356-27-6	59
4			Terephthalic acid, di(4-octyl) e...	390	C24H38O4	1000323-74-2	49
5			1,3-Benzenedicarboxylic acid, bi...	390	C24H38O4	000137-89-3	49



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG102722\
Data File : BG055332.D
Acq On : 27 Oct 2022 20:46
Operator : CG/JU
Sample : N5282-01
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SUMP-STONE

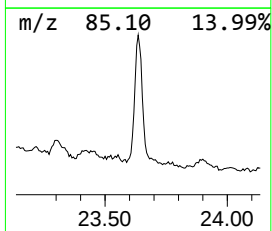
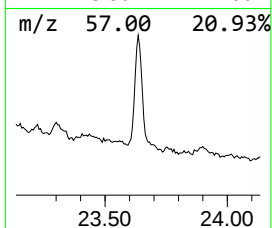
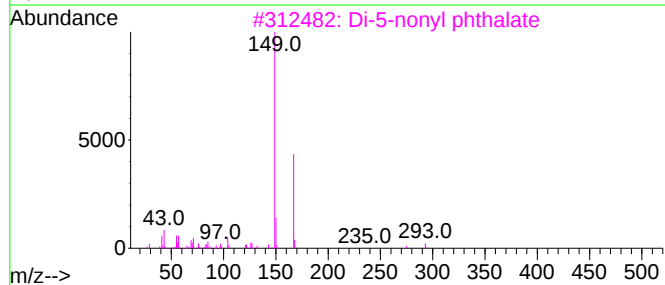
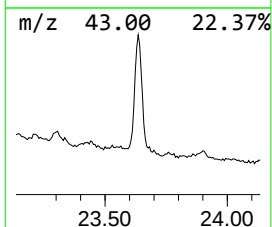
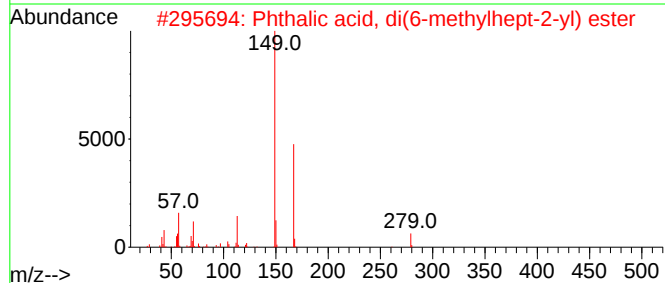
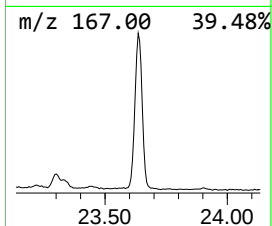
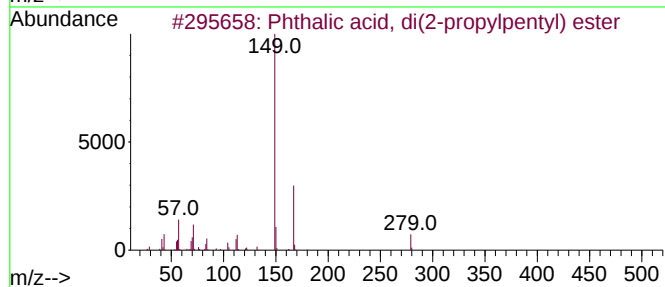
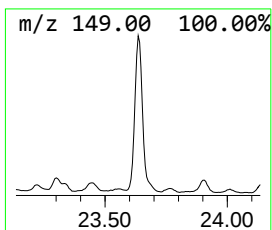
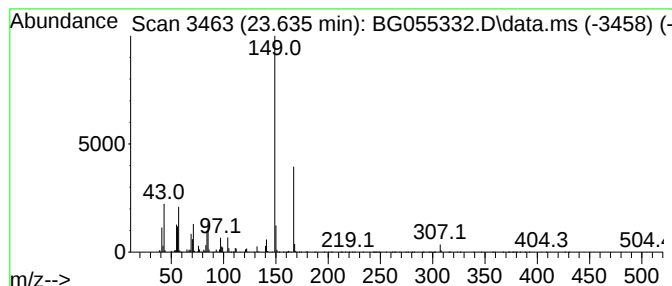
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG102422.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Phthalic acid, di(2-propylp... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.635	20.00 ng	1540020	Perylene-d12	25.321

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phthalic acid, di(2-propylpentyl...	390	C24H38O4	1000377-93-5	80
2			Phthalic acid, di(6-methylhept-2...	390	C24H38O4	1000377-97-3	72
3			Di-5-nonyl phthalate	418	C26H42O4	094054-23-6	72
4			Phthalic acid, monocyclohexyl ester	248	C14H16O4	007517-36-4	64
5			Bis(2-ethylhexyl) phthalate	390	C24H38O4	000117-81-7	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG102722\
Data File : BG055332.D
Acq On : 27 Oct 2022 20:46
Operator : CG/JU
Sample : N5282-01
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SUMP-STONE

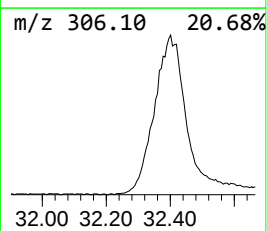
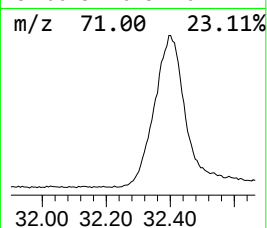
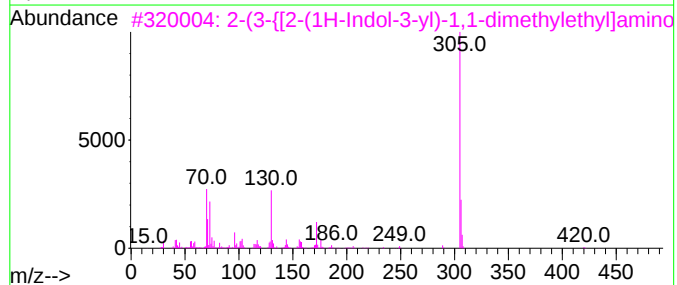
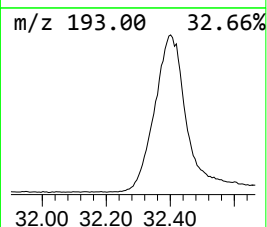
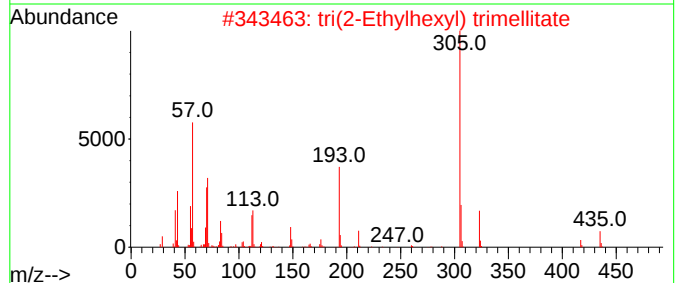
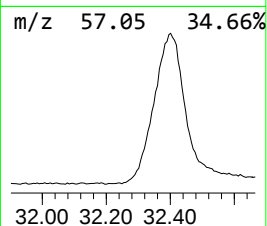
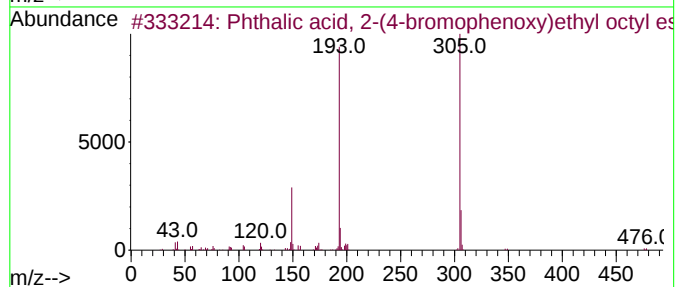
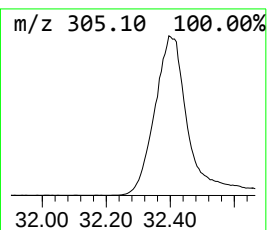
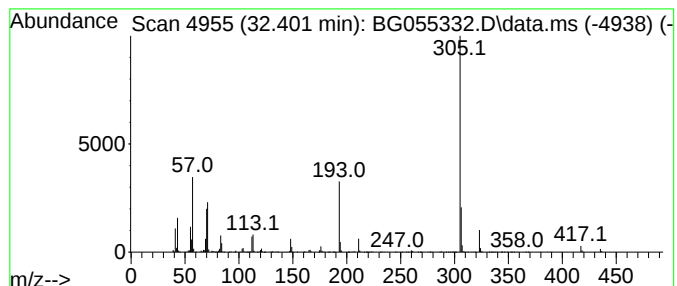
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG102422.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Phthalic acid, 2-(4-bromoph... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
32.401	52.27 ng	4024990	Perylene-d12	25.321

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phthalic acid, 2-(4-bromophenoxy...	476	C24H29BrO5	1010382-89-7	50
2			tri(2-Ethylhexyl) trimellitate	546	C33H54O6	003319-31-1	38
3			2-(3-{[2-(1H-Indol-3-yl)-1,1-dim...	435	C25H33N3O2Si	1000408-41-9	36
4			Terephthalic acid, octyl 2-pheno...	398	C24H30O5	1000416-04-7	28
5			4-(Diphenylphosphino)-N,N-dimeth...	305	C20H20NP	000739-58-2	25



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG102722\
Data File : BG055332.D
Acq On : 27 Oct 2022 20:46
Operator : CG/JU
Sample : N5282-01
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SUMP-STONE

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG102422.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST0.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-...	5.315	8.8	ng	129554	1	8.271	293721	20.0
1,3,4-Trimethyl...	12.184	5.3	ng	107327	2	11.103	407947	20.0
1,1'-Bicyclohexyl	12.871	5.5	ng	112871	2	11.103	407947	20.0
Naphthalene, 1,...	12.977	10.0	ng	204275	2	11.103	407947	20.0
Dodecane, 2,6,1...	13.406	3.8	ng	311351	3	14.893	1630640	20.0
unknown13.676	13.676	7.1	ng	580945	3	14.893	1630640	20.0
1H-Indene, octa...	14.311	14.4	ng	1176250	3	14.893	1630640	20.0
Tetradecane	14.346	14.5	ng	1183590	3	14.893	1630640	20.0
3-tert-Butyl-4-...	14.399	9.5	ng	772191	3	14.893	1630640	20.0
2-Dodecen-1-yl(...	14.634	18.3	ng	1491620	3	14.893	1630640	20.0
unknown14.722	14.722	20.0	ng	1630640	3	14.893	1630640	20.0
1,4-Benzenedica...	23.095	20.0	ng	590053	5	21.937	590053	20.0
Phthalic acid, ...	23.635	20.0	ng	1540020	6	25.321	1540020	20.0
Phthalic acid, ...	32.401	52.3	ng	4024990	6	25.321	1540020	20.0