

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020822\
 Data File : BP009059.D
 Acq On : 08 Feb 2022 15:31
 Operator : CG/JU
 Sample : N1421-03
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SB-3

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-BP020722.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BP009059.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.017	3	5	12	rVB	264392	262927	4.70%	0.709%
2	3.122	19	23	32	rVB	250716	297511	5.31%	0.802%
3	4.934	326	331	339	rBV	52582	80696	1.44%	0.218%
4	5.393	403	409	436	rBV	1948974	3102678	55.43%	8.365%
5	6.987	673	680	702	rBV	1720750	3148714	56.25%	8.489%
6	7.805	810	819	827	rBV	598450	1009147	18.03%	2.721%
7	8.963	1009	1016	1033	rBV	1225258	2256181	40.31%	6.083%
8	10.604	1286	1295	1312	rBV	799796	1436954	25.67%	3.874%
9	13.069	1706	1714	1736	rBV	3497213	5329782	95.22%	14.369%
10	14.445	1942	1948	1967	rVB	1077683	1714599	30.63%	4.622%
11	15.945	2196	2203	2222	rBV	2023763	3254687	58.14%	8.774%
12	17.204	2411	2417	2433	rBV	1073734	1734855	30.99%	4.677%
13	18.104	2565	2570	2582	rBV	162181	314179	5.61%	0.847%
14	19.339	2776	2780	2792	rBV3	93558	203339	3.63%	0.548%
15	19.769	2848	2853	2861	rBV	4682857	5597618	100.00%	15.091%
16	21.016	3062	3065	3072	rVB2	137090	196257	3.51%	0.529%
17	21.304	3109	3114	3121	rBV	1422602	1987027	35.50%	5.357%
18	21.416	3128	3133	3142	rBV	1910350	2652578	47.39%	7.151%
19	23.704	3516	3522	3537	rVB	1154731	2512925	44.89%	6.775%

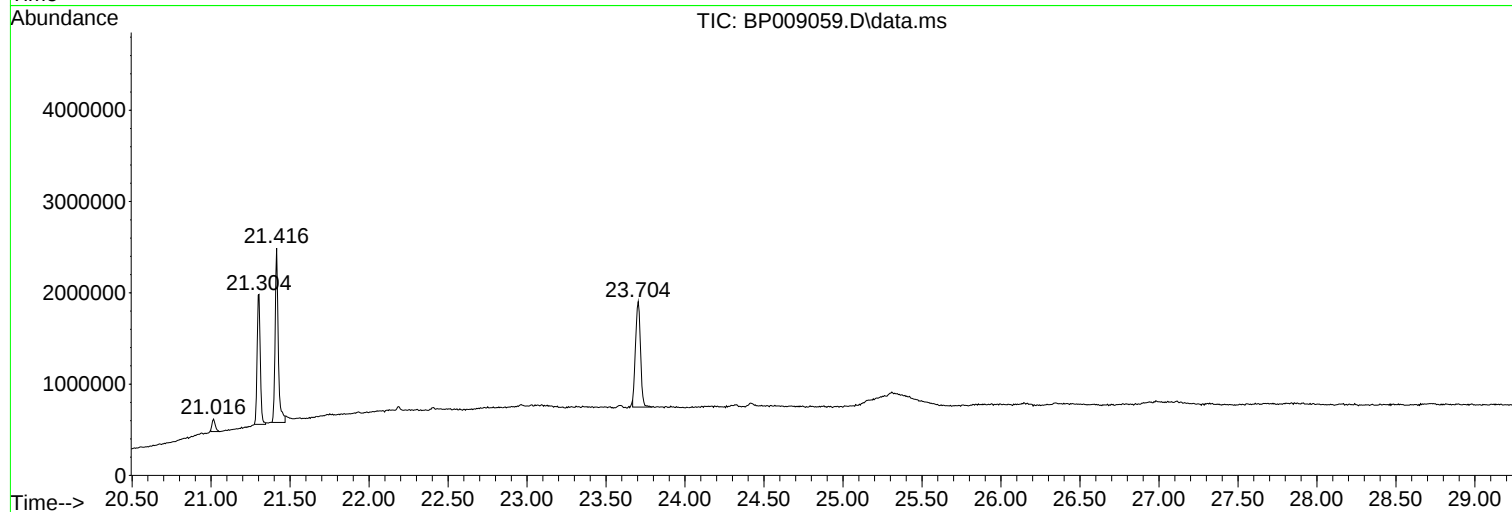
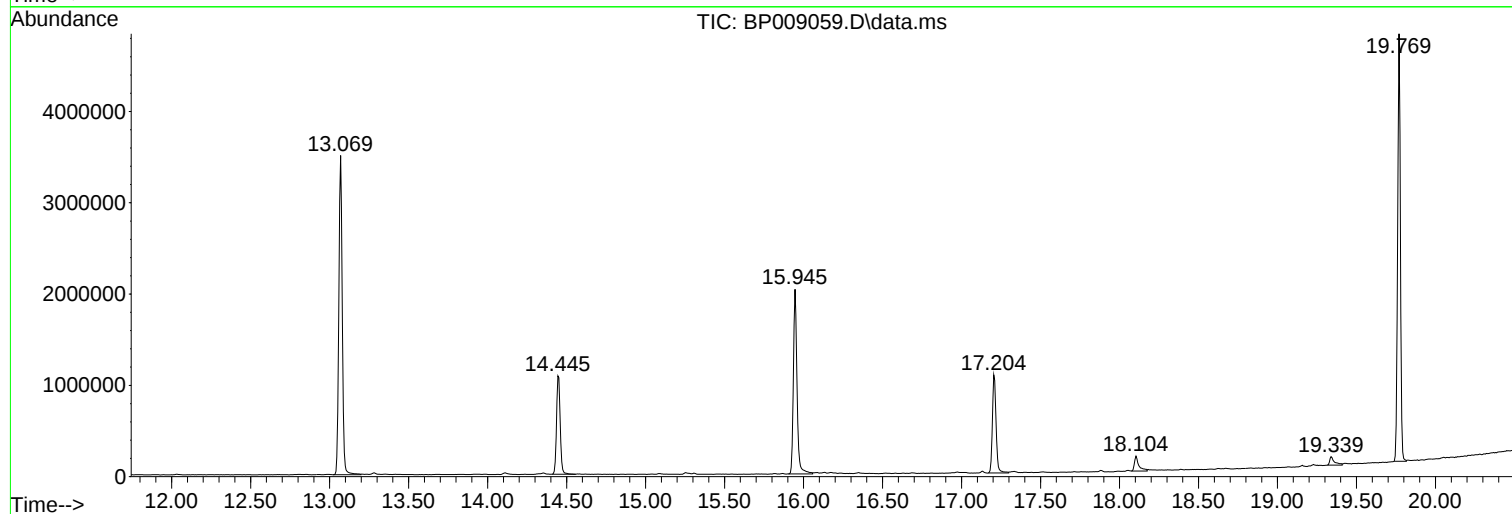
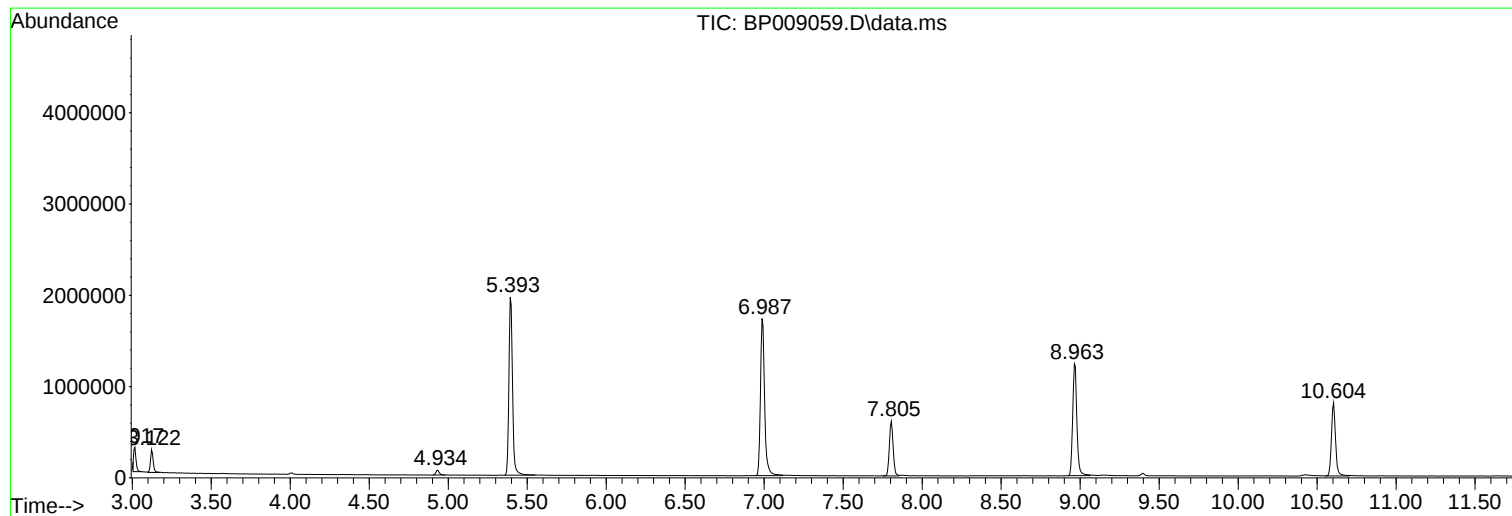
Sum of corrected areas: 37092654

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020822\
Data File : BP009059.D
Acq On : 08 Feb 2022 15:31
Operator : CG/JU
Sample : N1421-03
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
SB-3

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP020722.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\BP020822\
Data File : BP009059.D
Acq On : 08 Feb 2022 15:31
Operator : CG/JU
Sample : N1421-03
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
SB-3

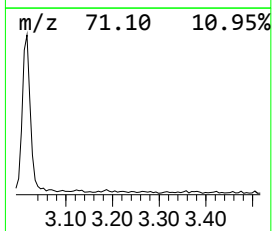
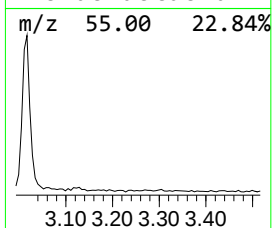
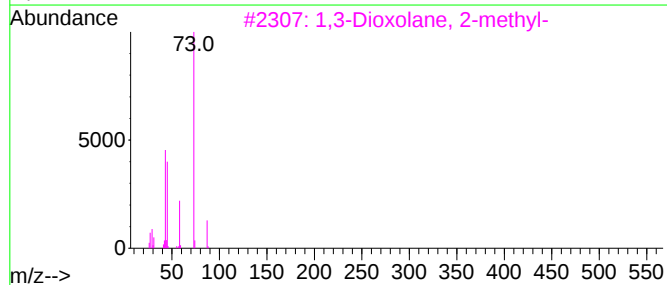
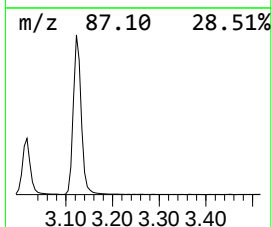
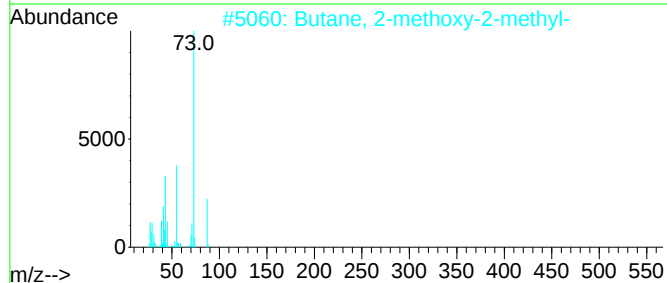
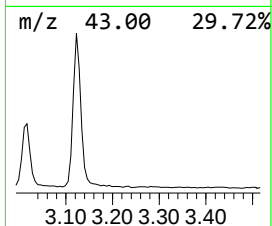
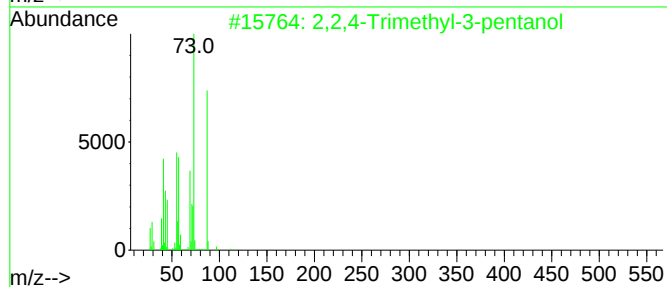
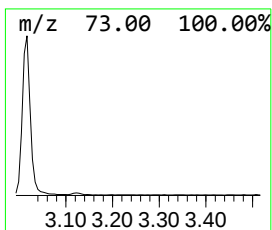
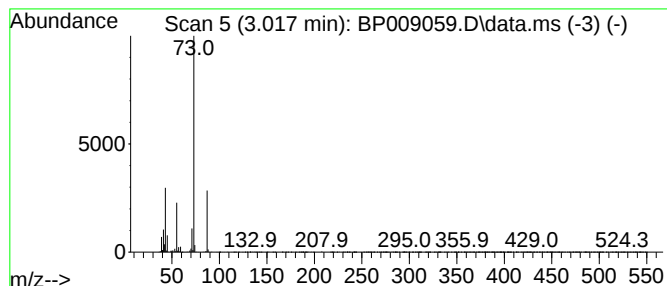
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP020722.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 unknown3.017 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.017	5.21 ng	262927	1,4-Dichlorobenzene-d4	7.805

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	40
2			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	40
3			1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	9
4			Silane, tetramethyl-	88	C4H12Si	000075-76-3	9
5			N-Ethylformamide	73	C3H7NO	000627-45-2	9



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020822\
Data File : BP009059.D
Acq On : 08 Feb 2022 15:31
Operator : CG/JU
Sample : N1421-03
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
SB-3

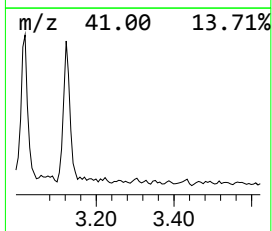
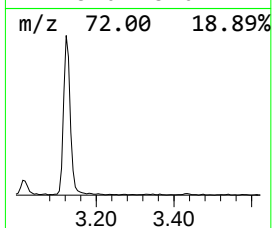
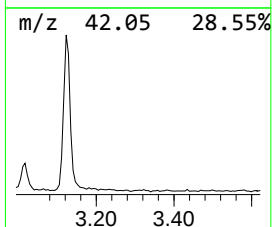
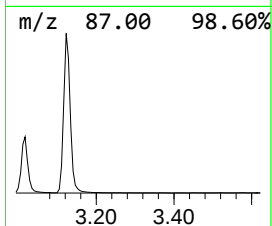
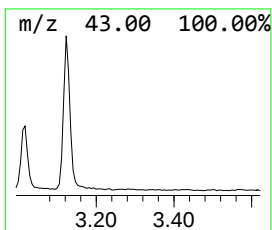
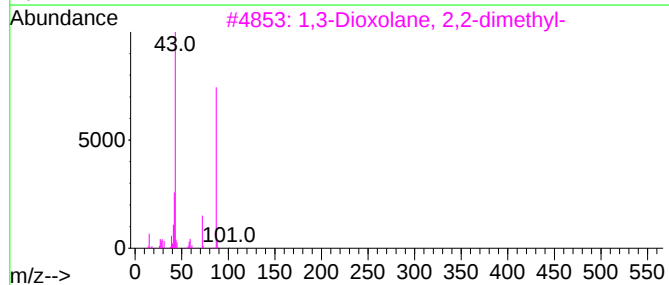
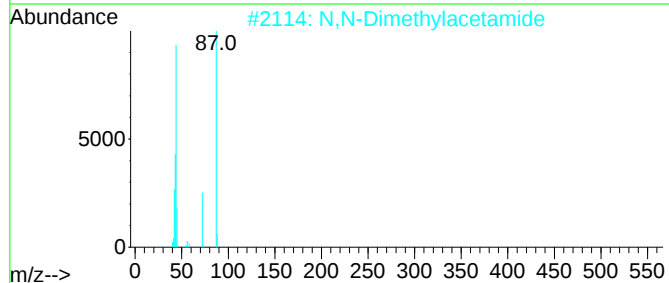
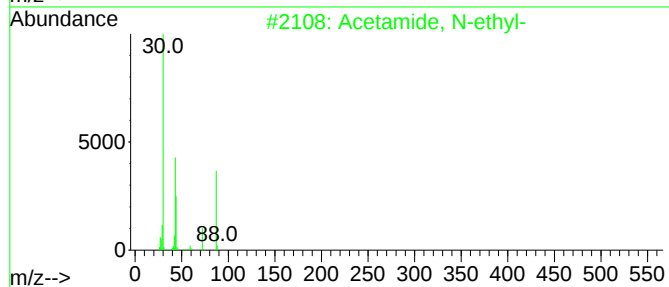
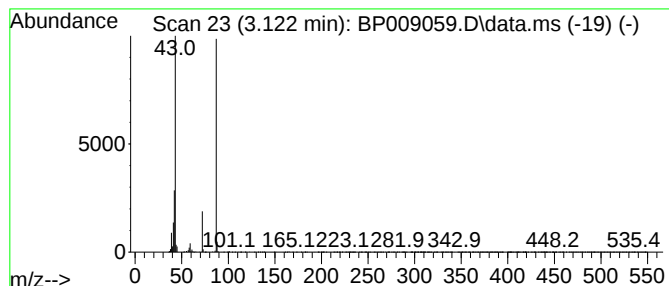
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP020722.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 Acetamide, N-ethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.122	5.90 ng	297511	1,4-Dichlorobenzene-d4	7.805

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	72
2			N,N-Dimethylacetamide	87	C4H9NO	000127-19-5	72
3			1,3-Dioxolane, 2,2-dimethyl-	102	C5H10O2	002916-31-6	53
4			2-OXOPROPIONAMIDE	87	C3H5NO2	000631-66-3	50
5			1,3-Dioxolane-2-ethanol, 2-methyl-	132	C6H12O3	005754-32-5	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020822\
Data File : BP009059.D
Acq On : 08 Feb 2022 15:31
Operator : CG/JU
Sample : N1421-03
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
SB-3

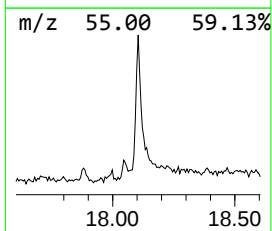
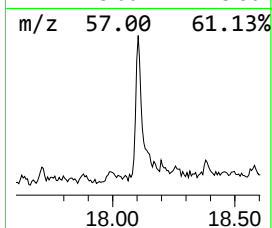
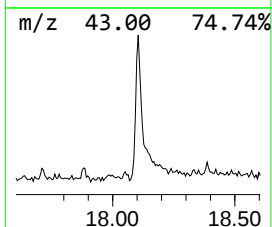
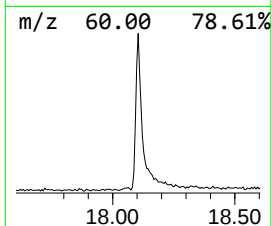
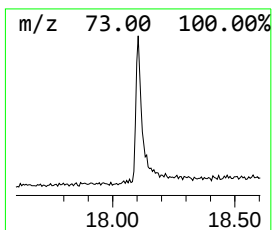
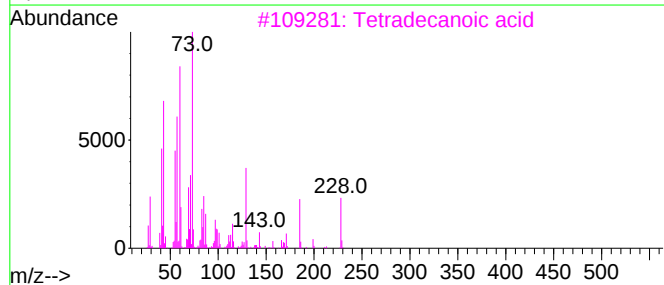
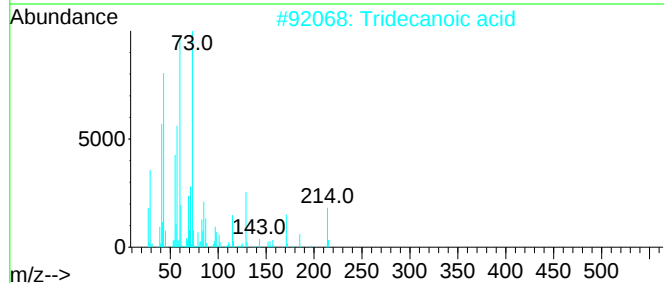
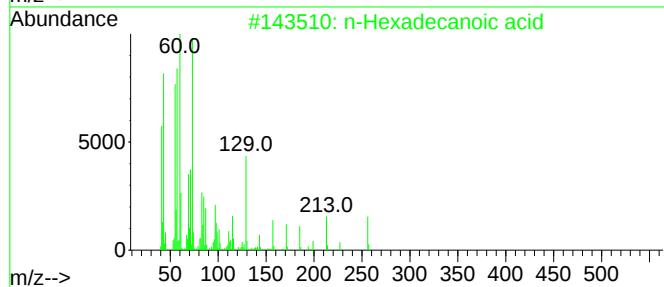
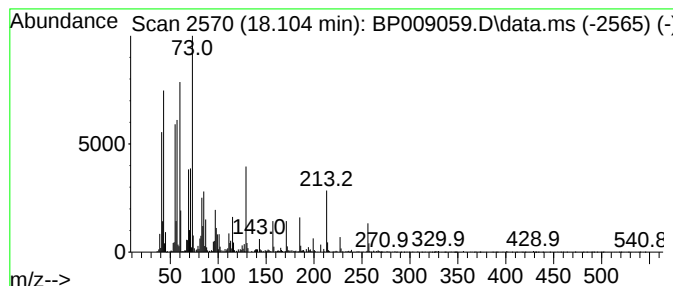
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP020722.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 3 n-Hexadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.104	3.62 ng	314179	Phenanthrene-d10	17.204

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2		Tridecanoic acid	214	C13H26O2	000638-53-9	93
3		Tetradecanoic acid	228	C14H28O2	000544-63-8	93
4		Octadecanoic acid	284	C18H36O2	000057-11-4	87
5		n-Decanoic acid	172	C10H20O2	000334-48-5	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020822\
Data File : BP009059.D
Acq On : 08 Feb 2022 15:31
Operator : CG/JU
Sample : N1421-03
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
SB-3

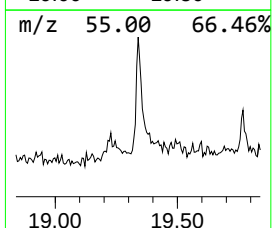
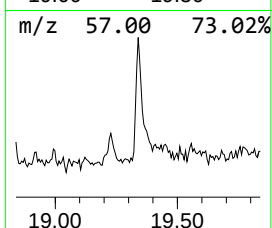
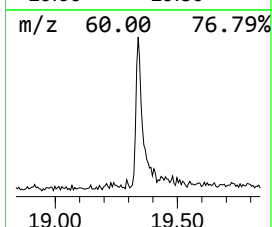
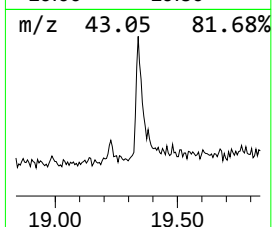
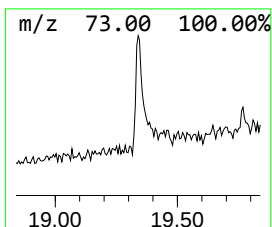
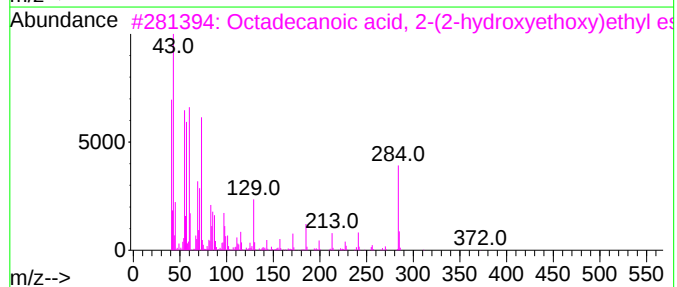
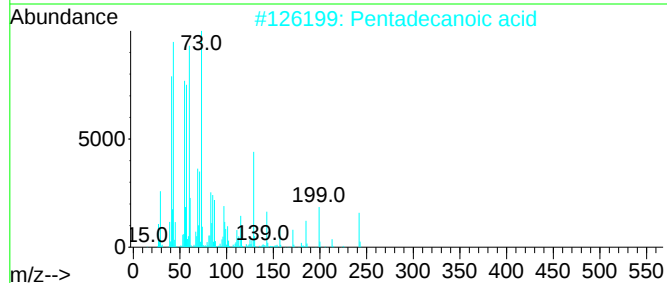
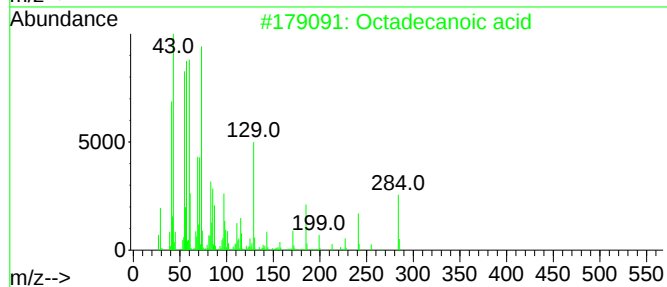
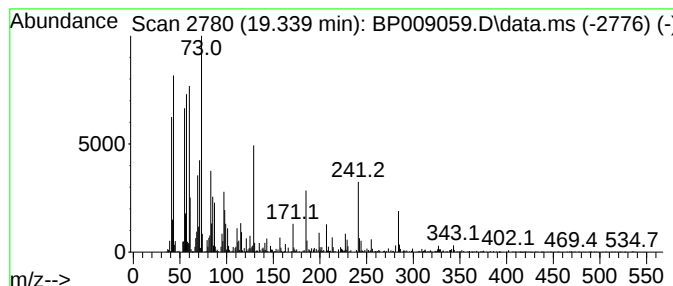
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP020722.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 4 Octadecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.339	2.05 ng	203339	Chrysene-d12	21.304

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid	284	C18H36O2	000057-11-4	99
2			Pentadecanoic acid	242	C15H30O2	001002-84-2	90
3			Octadecanoic acid, 2-(2-hydroxye...	372	C22H44O4	000106-11-6	86
4			Docosanoic acid	340	C22H44O2	000112-85-6	83
5			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020822\
Data File : BP009059.D
Acq On : 08 Feb 2022 15:31
Operator : CG/JU
Sample : N1421-03
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
SB-3

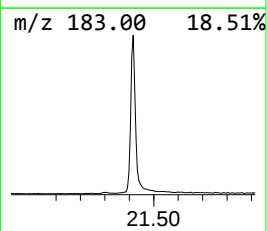
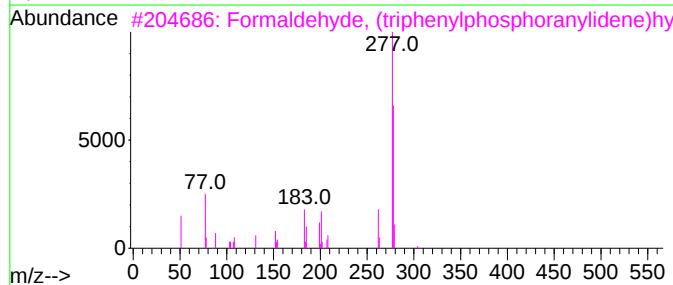
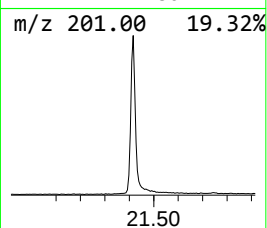
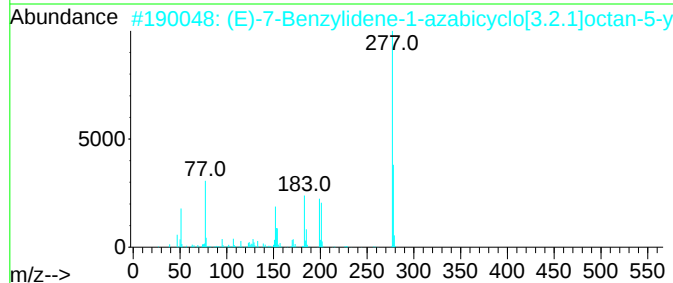
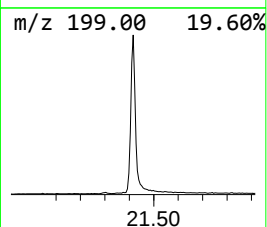
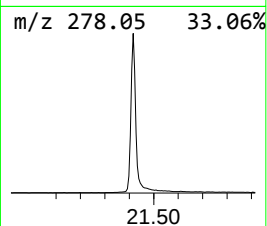
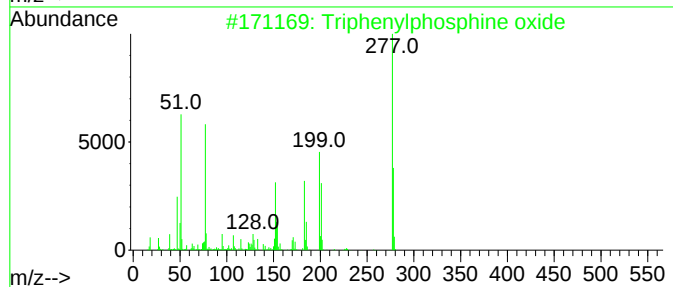
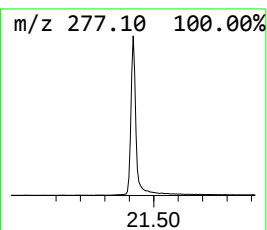
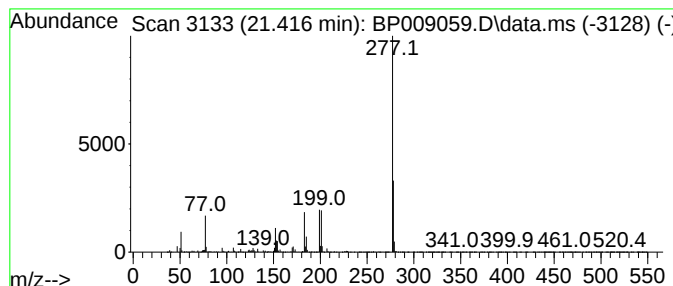
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP020722.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Triphenylphosphine oxide Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.416	26.70 ng	2652580	Chrysene-d12	21.304

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Triphenylphosphine oxide	278	C18H15OP	000791-28-6	95
2			(E)-7-Benzylidene-1-azabicyclo[3...]	293	C15H19NO3S	1000483-83-4	90
3			Formaldehyde, (triphenylphosphor...	304	C19H17N2P	015990-54-2	45
4			2-Phenyl-6-(trifluoromethyl)-1H-...	277	C15H10F3NO	1000387-59-3	43
5			3-(3,4-Dichlorophenyl)-2-thioxo-...	277	C9H5Cl2NOS2	017046-34-3	32



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020822\
Data File : BP009059.D
Acq On : 08 Feb 2022 15:31
Operator : CG/JU
Sample : N1421-03
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
SB-3

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP020722.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
unknown3.017	3.017	5.2	ng	262927	1	7.805	1009150	20.0
Acetamide, N-et...	3.122	5.9	ng	297511	1	7.805	1009150	20.0
n-Hexadecanoic ...	18.104	3.6	ng	314179	4	17.204	1734860	20.0
Octadecanoic acid	19.339	2.0	ng	203339	5	21.304	1987030	20.0
Triphenylphosph...	21.416	26.7	ng	2652580	5	21.304	1987030	20.0