

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP022122\
 Data File : BP009231.D
 Acq On : 21 Feb 2022 13:13
 Operator : CG/JU
 Sample : N1599-01
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SPN-4

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: BP009231.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.246	209	214	232	rVB	159375	291875	2.75%	0.658%
2	5.363	398	404	426	rBV	661131	1337565	12.60%	3.017%
3	6.969	671	677	695	rBV	352837	877040	8.26%	1.978%
4	7.316	730	736	741	rBV	90453	150189	1.41%	0.339%
5	7.369	741	745	753	rVB	91718	149699	1.41%	0.338%
6	7.557	770	777	789	rBV	132317	268052	2.52%	0.605%
7	7.763	805	812	818	rBV	300455	537774	5.06%	1.213%
8	7.828	818	823	840	rVB	136674	271629	2.56%	0.613%
9	8.928	1003	1010	1034	rBV	867713	1753888	16.52%	3.956%
10	10.557	1280	1287	1303	rBV	383289	765127	7.21%	1.726%
11	12.763	1653	1662	1676	rVB	115563	227612	2.14%	0.513%
12	13.028	1698	1707	1726	rBV	3174949	5324816	50.15%	12.012%
13	13.263	1738	1747	1762	rBV	655321	1122864	10.57%	2.533%
14	14.410	1935	1942	1952	rBV	741581	1211938	11.41%	2.734%
15	15.233	2074	2082	2090	rVB	134668	218715	2.06%	0.493%
16	15.780	2169	2175	2185	rBV2	151177	297791	2.80%	0.672%
17	15.916	2191	2198	2213	rBV	2498419	4387969	41.33%	9.898%
18	17.174	2405	2412	2425	rBV	950505	1577455	14.86%	3.558%
19	17.833	2519	2524	2530	rBV2	171601	220882	2.08%	0.498%
20	18.074	2559	2565	2588	rBV	1919808	3424844	32.25%	7.726%
21	19.192	2747	2755	2766	rBV7	110558	399515	3.76%	0.901%
22	19.304	2769	2774	2797	rVV	1214637	2281742	21.49%	5.147%
23	19.733	2841	2847	2859	rBV	8318006	10618103	100.00%	23.952%
24	20.968	3054	3057	3066	rBV2	156314	226736	2.14%	0.511%
25	21.045	3067	3070	3076	rVV	92658	119789	1.13%	0.270%
26	21.168	3088	3091	3096	rVB	219335	232891	2.19%	0.525%
27	21.268	3104	3108	3117	rBV	1646103	2245255	21.15%	5.065%
28	22.486	3311	3315	3322	rVB	546091	839491	7.91%	1.894%
29	23.651	3507	3513	3527	rVB2	1329859	2948966	27.77%	6.652%

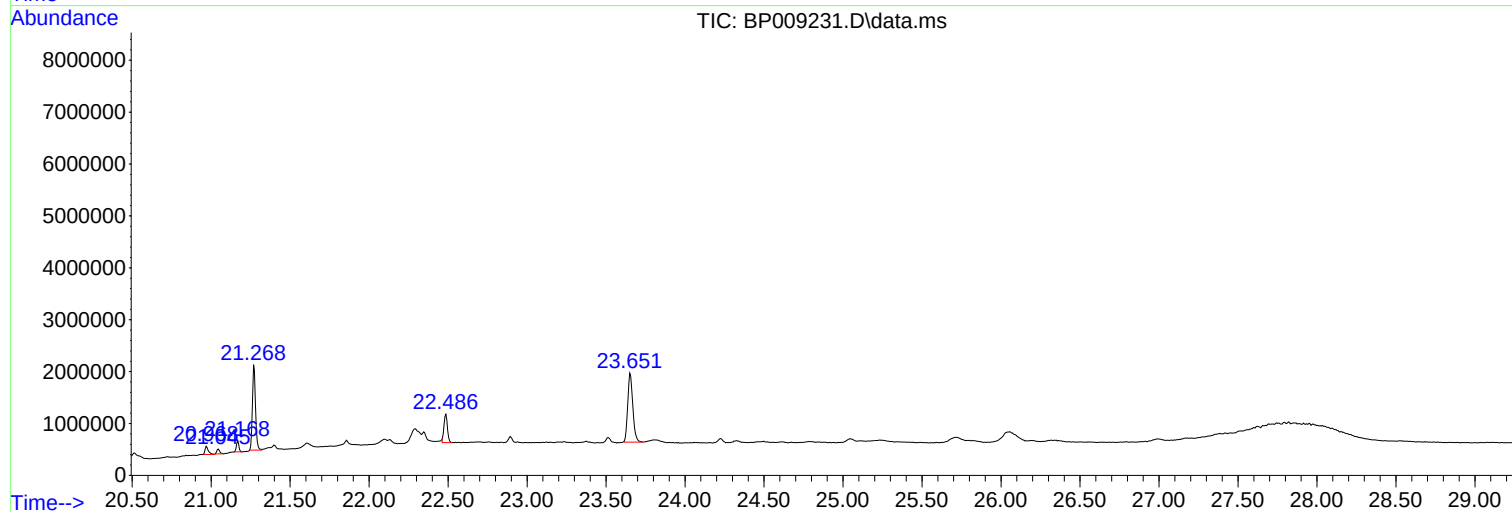
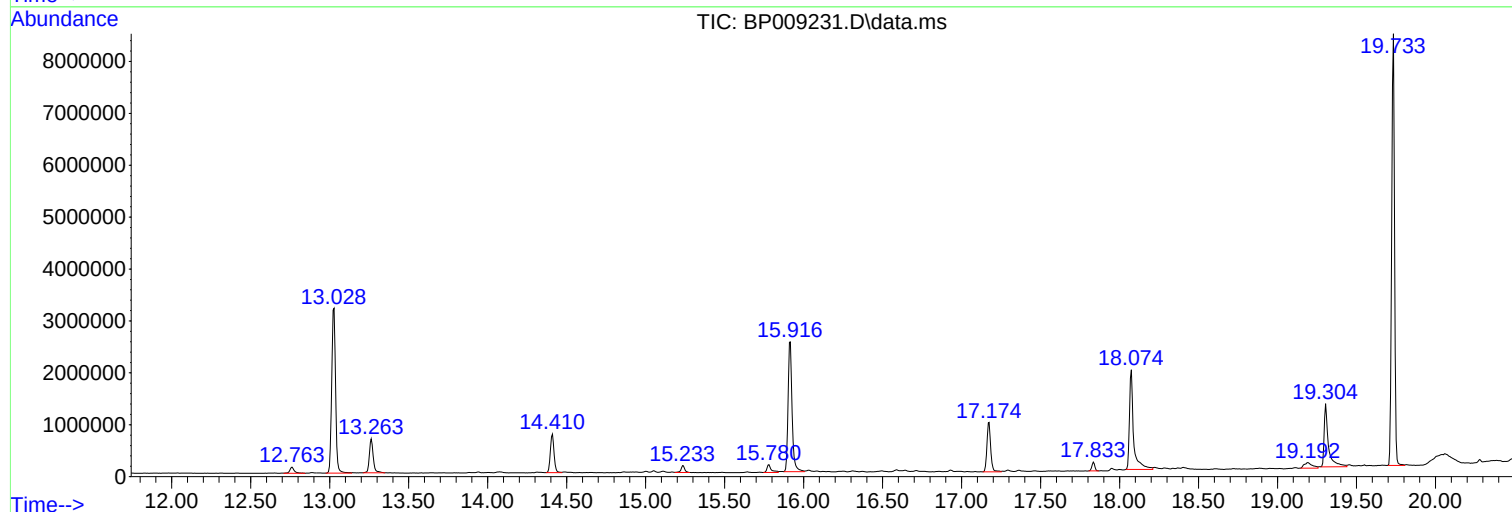
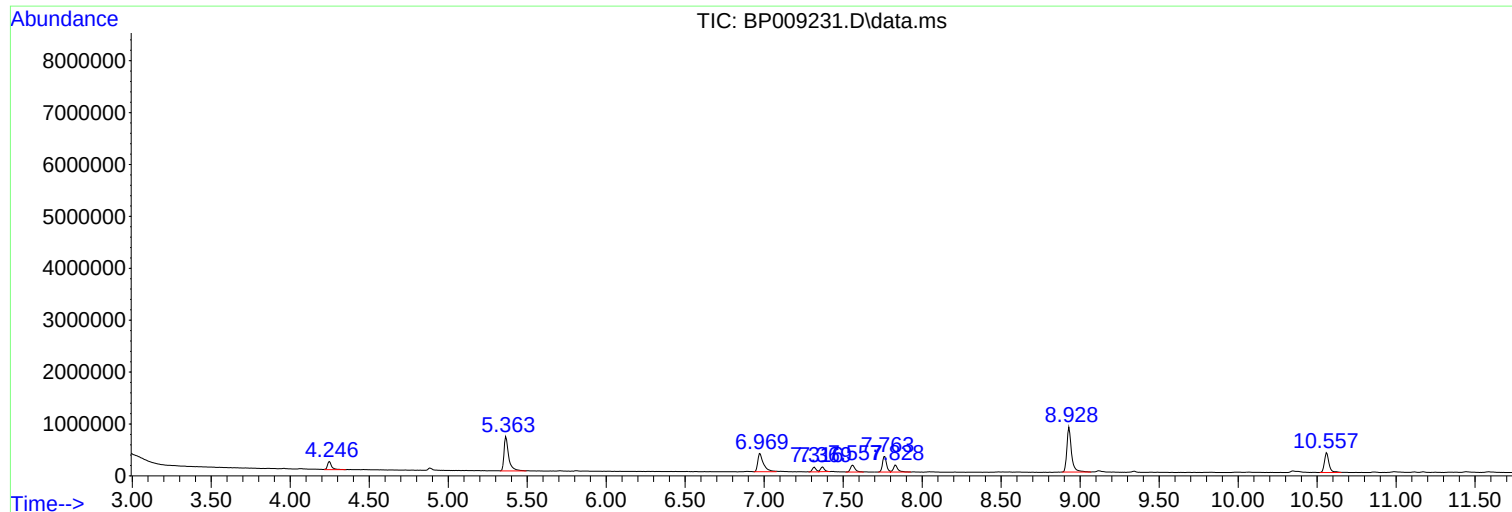
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BNA_P
ClientSampleId :
SPN-4

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



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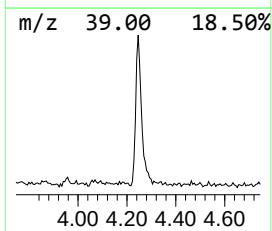
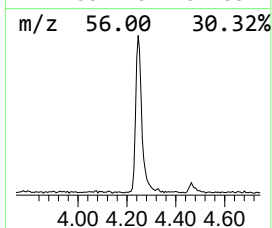
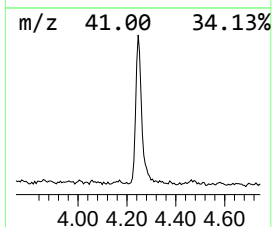
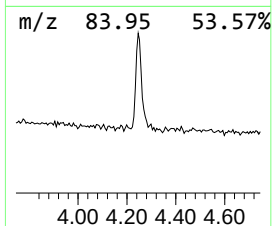
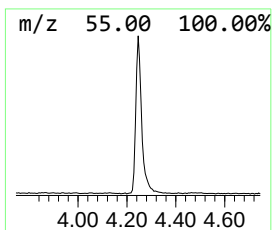
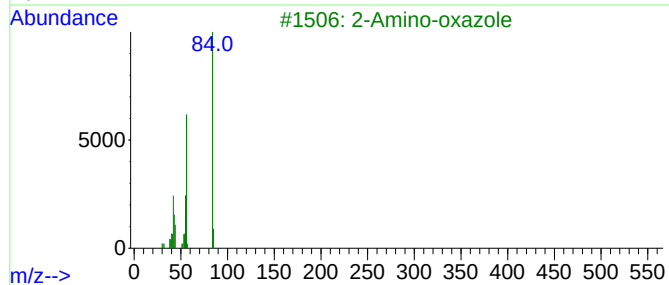
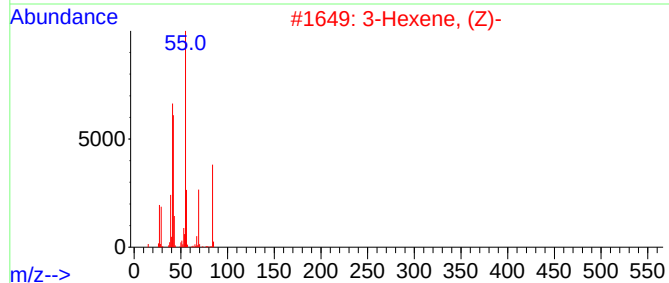
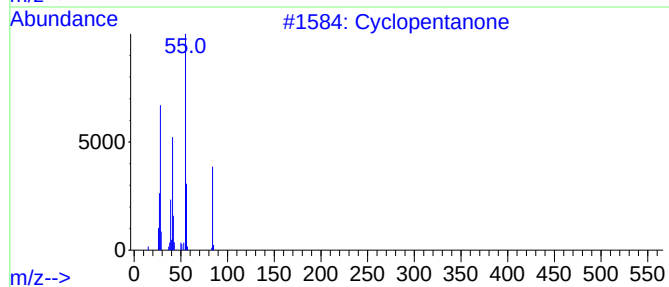
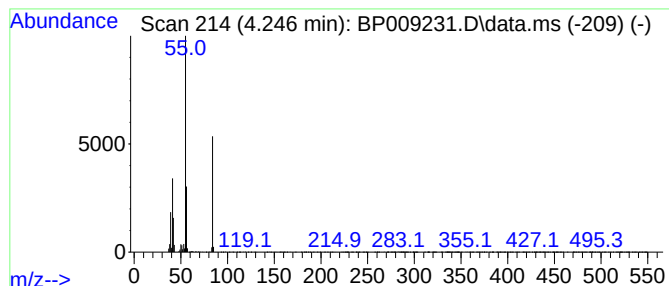
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Cyclopentanone Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.246	10.85 ng	291875	1,4-Dichlorobenzene-d4	7.763

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopentanone	84	C5H8O	000120-92-3	91
2			3-Hexene, (Z)-	84	C6H12	007642-09-3	53
3			2-Amino-oxazole	84	C3H4N2O	004570-45-0	53
4			2(5H)-Furanone	84	C4H4O2	000497-23-4	50
5			2H-Pyran, 3,4-dihydro-	84	C5H8O	000110-87-2	45



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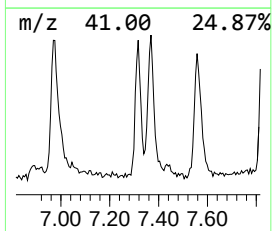
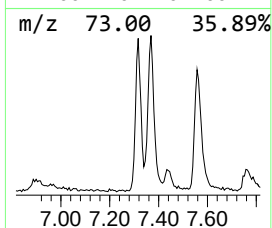
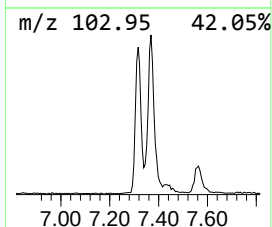
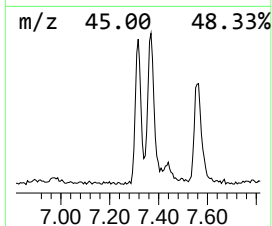
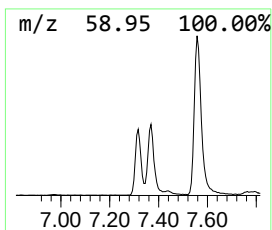
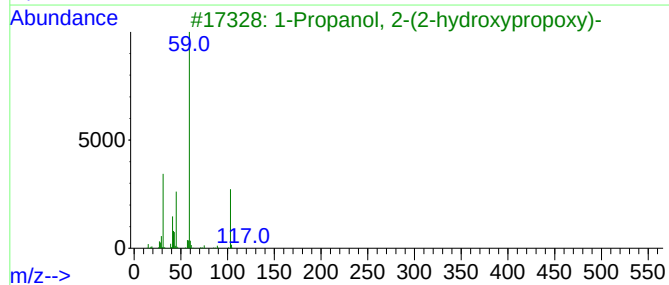
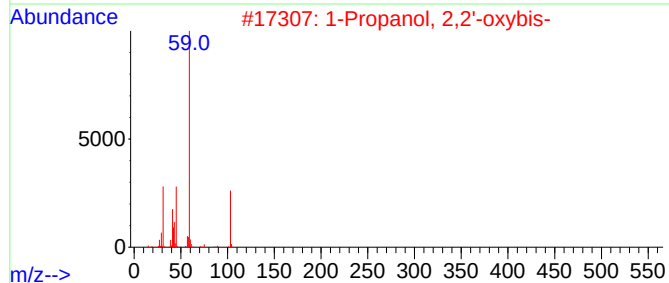
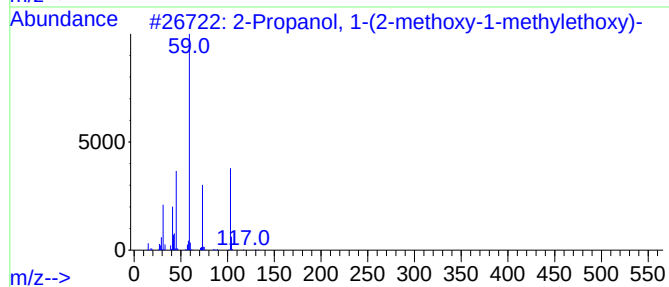
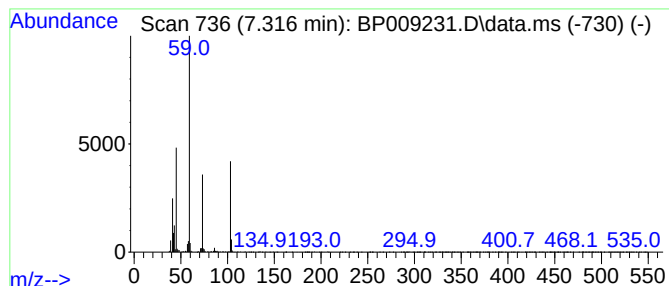
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP020722.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 2-Propanol, 1-(2-methoxy-1-... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.316	5.59 ng	150189	1,4-Dichlorobenzene-d4	7.763

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Propanol, 1-(2-methoxy-1-methy...	148	C7H16O3	020324-32-7	90
2			1-Propanol, 2,2'-oxybis-	134	C6H14O3	000108-61-2	50
3			1-Propanol, 2-(2-hydroxypropoxy)-	134	C6H14O3	000106-62-7	50
4			Dipropylene glycol (isomer 3)	134	C6H14O3	1000506-25-5	50
5			Ethyl 2-(2-(2-methoxyethoxy)etho...	206	C9H18O5	1000366-89-0	40



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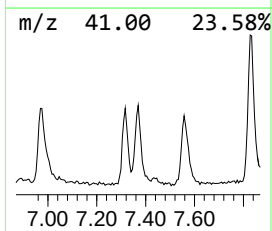
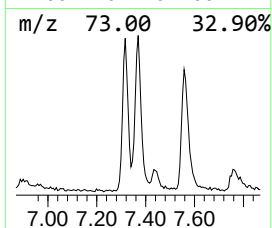
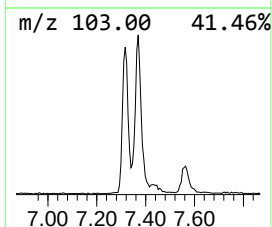
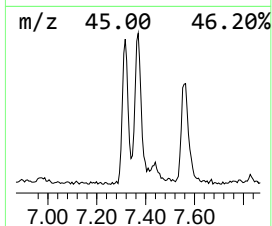
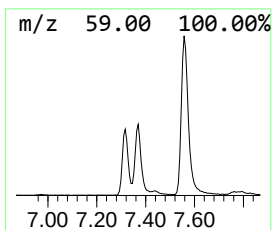
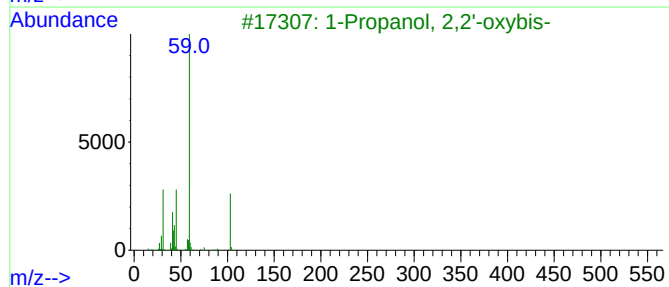
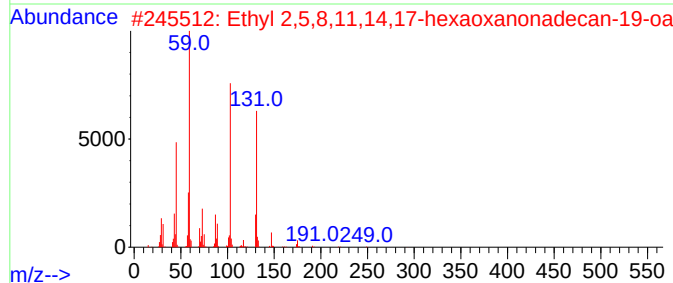
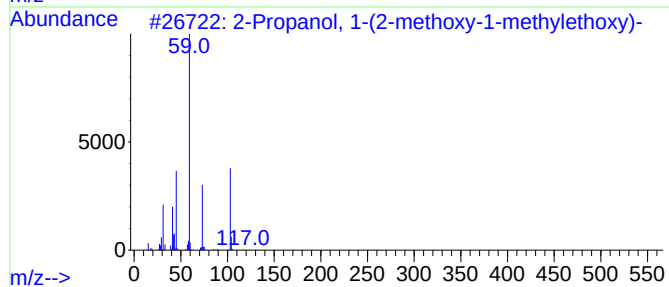
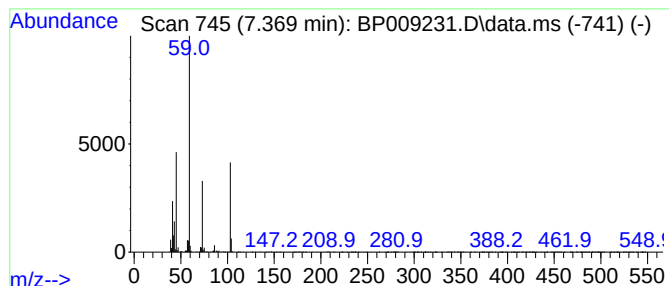
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Ethyl 2,5,8,11,14,17-hexaox... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.369	5.57 ng	149699	1,4-Dichlorobenzene-d4	7.763

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Propanol, 1-(2-methoxy-1-methy...	148	C7H16O3	020324-32-7	83		
2	Ethyl 2,5,8,11,14,17-hexaoxanona...	338	C15H30O8	1000366-80-6	56		
3	1-Propanol, 2,2'-oxybis-	134	C6H14O3	000108-61-2	53		
4	Dipropylene glycol (isomer 3)	134	C6H14O3	1000506-25-5	53		
5	Dipropylene glycol (isomer 2)	134	C6H14O3	1000506-25-4	53		



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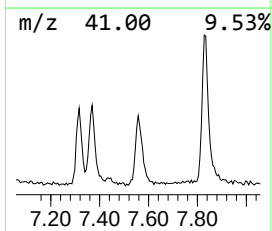
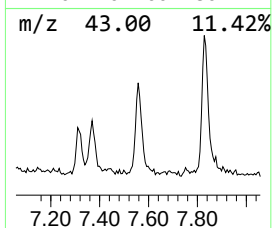
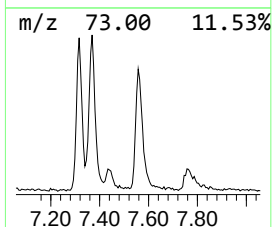
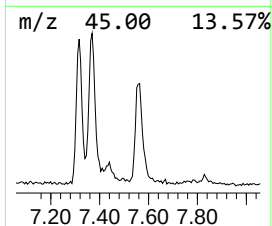
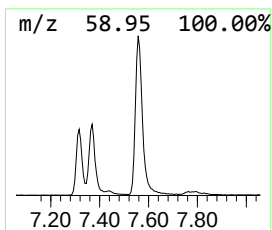
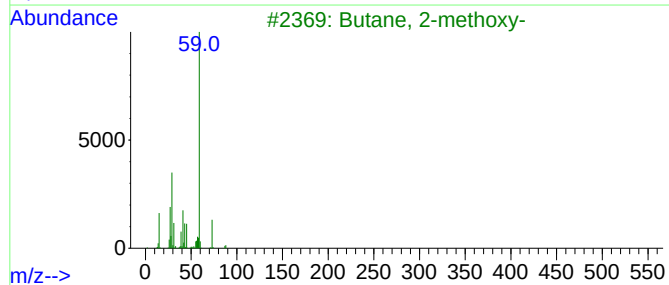
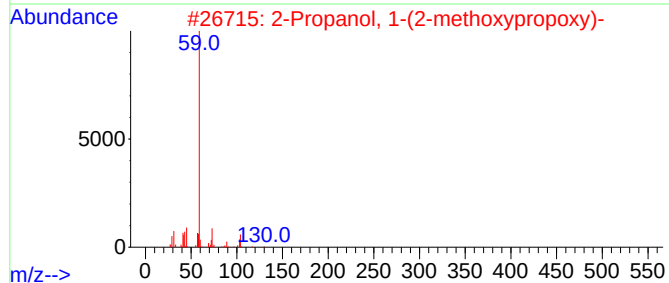
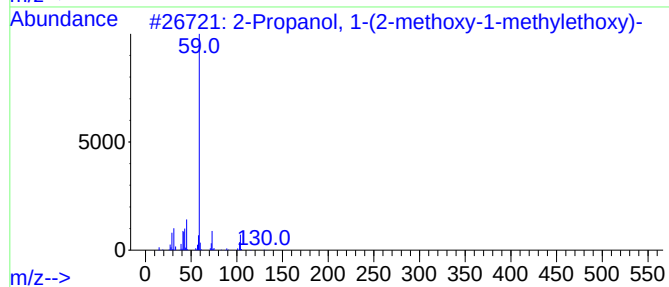
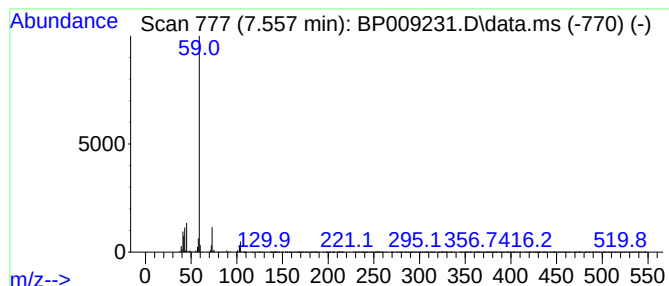
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Peak Number 4 2-Propanol, 1-(2-methoxypro... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD			R.T.
7.557	9.97 ng	268052	1,4-Dichlorobenzene-d4			7.763

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Propanol, 1-(2-methoxy-1-methy...	148	C7H16O3		020324-32-7	83
2	2-Propanol, 1-(2-methoxypropoxy)-	148	C7H16O3		013429-07-7	83
3	Butane, 2-methoxy-	88	C5H12O		006795-87-5	40
4	1-(1-Methoxypropan-2-yloxy)propa...	190	C9H18O4		1000367-08-6	40
5	Butane, 1-ethoxy-	102	C6H14O		000628-81-9	39



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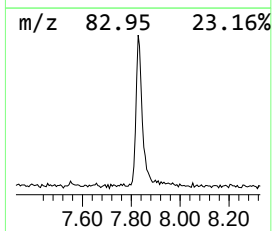
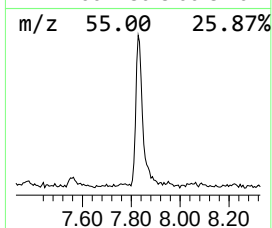
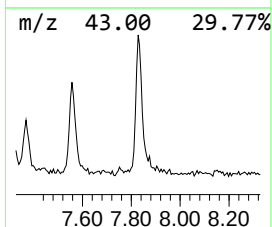
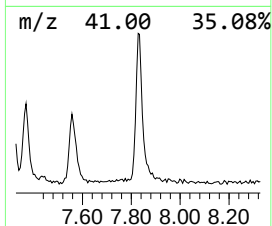
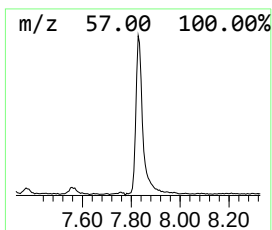
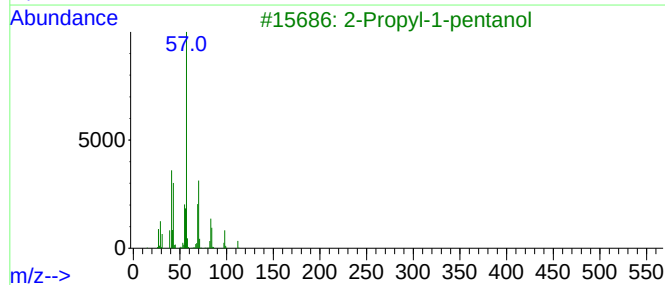
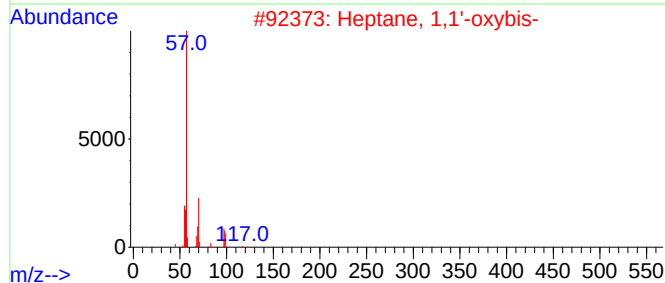
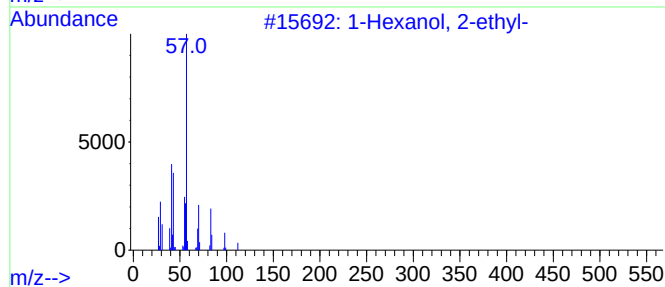
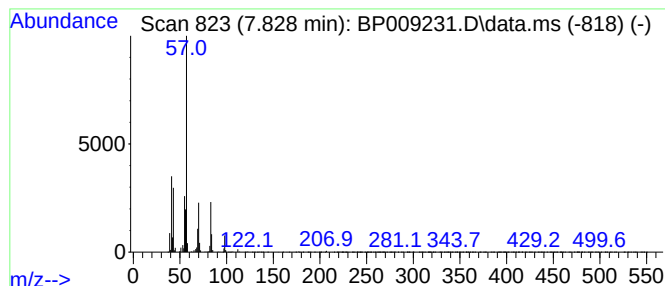
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 1-Hexanol, 2-ethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.828	10.10 ng	271629	1,4-Dichlorobenzene-d4	7.763

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	90
2		Heptane, 1,1'-oxybis-	214	C14H30O	000629-64-1	59
3		2-Propyl-1-pentanol	130	C8H18O	058175-57-8	50
4		2-Ethyl-1-hexanol, methyl ether	144	C9H20O	1000365-10-2	50
5		2-Decene, 5-methyl-, (Z)-	154	C11H22	074645-86-6	47



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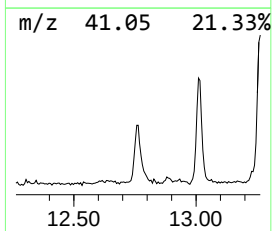
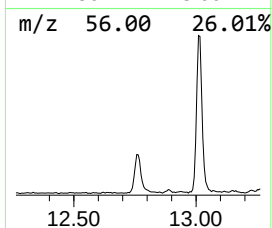
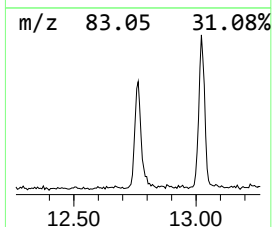
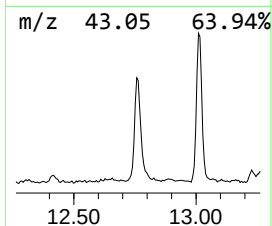
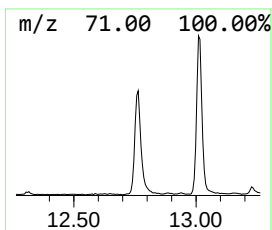
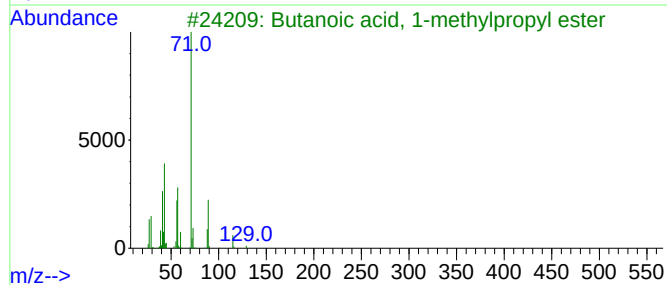
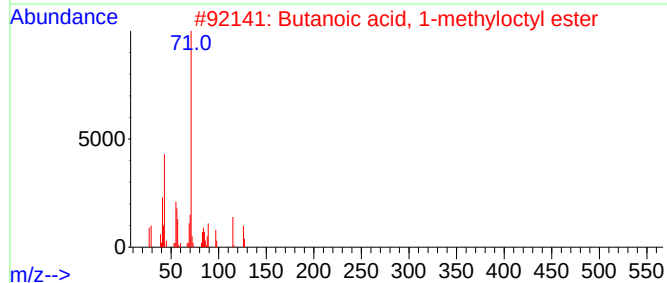
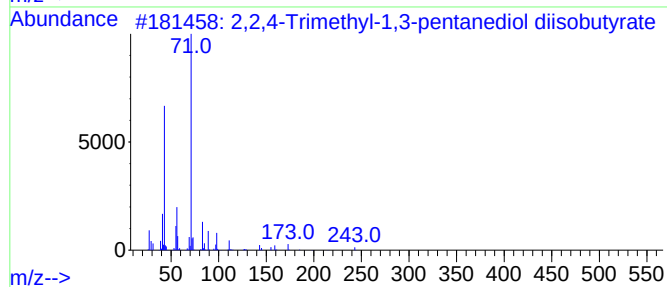
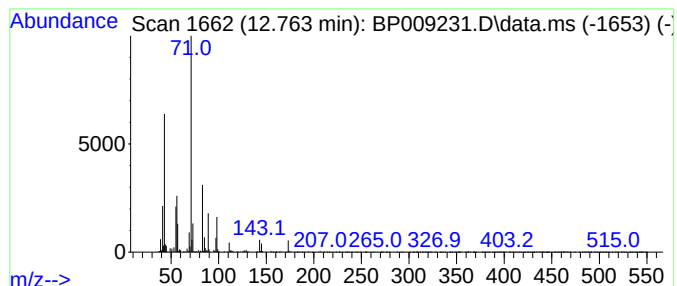
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 2,2,4-Trimethyl-1,3-pentane... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.763	3.76 ng	227612	Acenaphthene-d10	14.410

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2,2,4-Trimethyl-1,3-pentanediol ...	286	C16H30O4	006846-50-0	72
2			Butanoic acid, 1-methyloctyl ester	214	C13H26O2	069727-42-0	47
3			Butanoic acid, 1-methylpropyl ester	144	C8H16O2	000819-97-6	42
4			Isophytol	296	C20H40O	000505-32-8	38
5			Butyric acid, 3-tetradecyl ester	284	C18H36O2	1000280-57-1	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP022122\
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Acq On : 21 Feb 2022 13:13
Operator : CG/JU
Sample : N1599-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

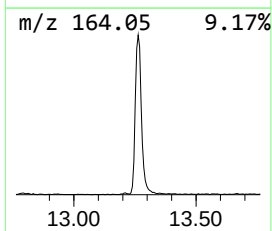
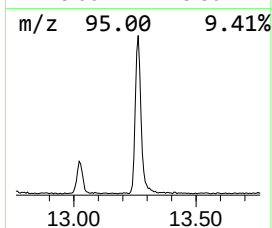
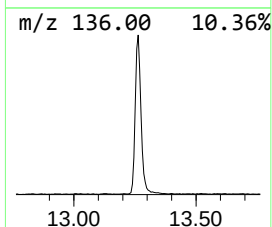
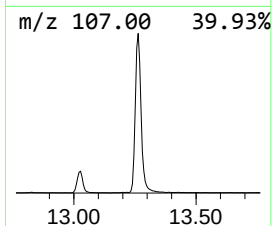
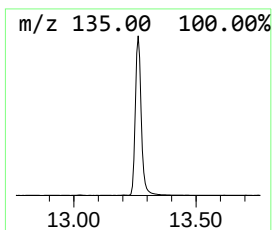
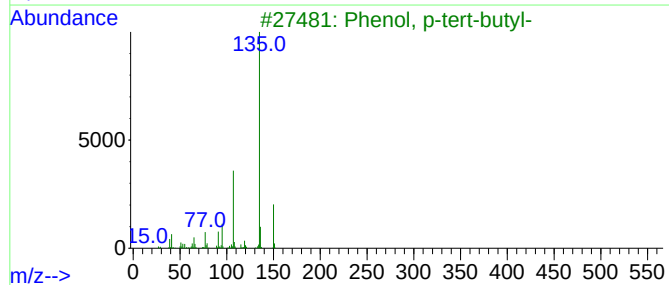
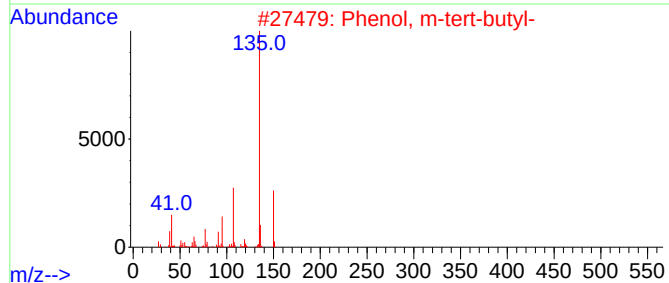
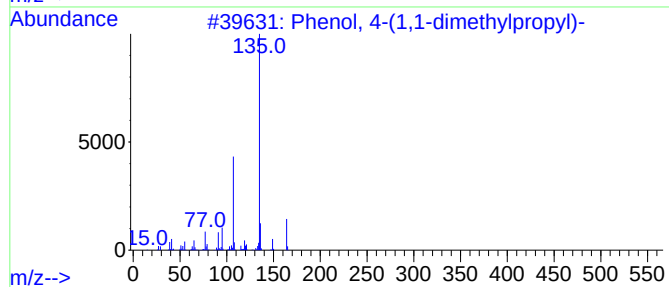
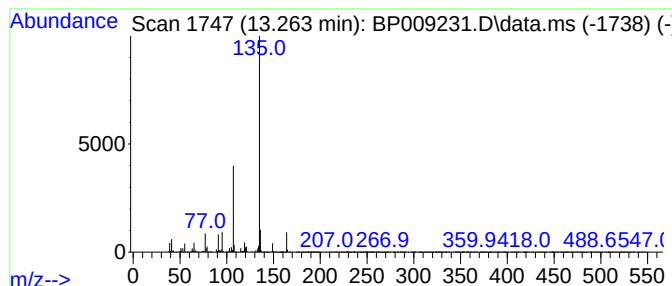
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Phenol, 4-(1,1-dimethylprop... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD		R.T.	
13.263	18.53 ng	1122860	Acenaphthene-d10		14.410	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Phenol, 4-(1,1-dimethylpropyl)-		164	C11H16O	000080-46-6	95
2	Phenol, m-tert-butyl-		150	C10H14O	000585-34-2	86
3	Phenol, p-tert-butyl-		150	C10H14O	000098-54-4	86
4	Imidazo[1,2-c]pyrimidin-5-ol		135	C6H5N3O	055662-66-3	64
5	4,6-Bis(4-ethoxybenzylthio)-5-ni...		457	C22H23N3O4S2	325957-76-4	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP022122\
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Sample : N1599-01
Misc :
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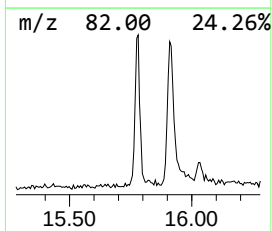
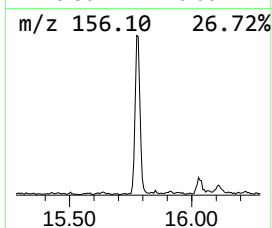
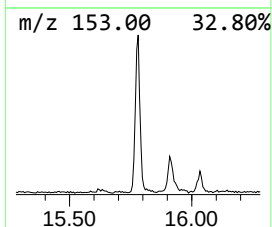
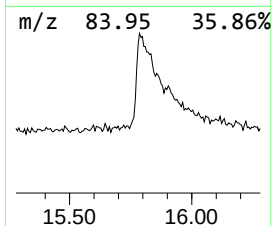
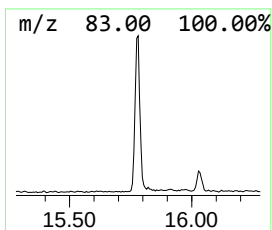
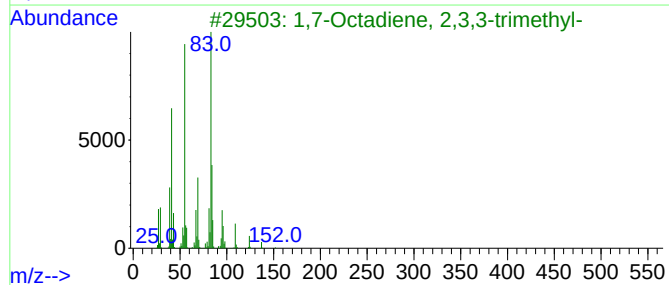
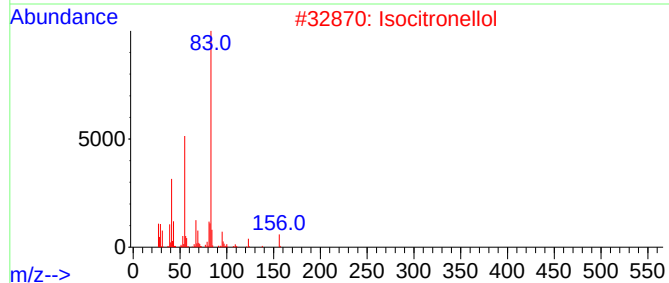
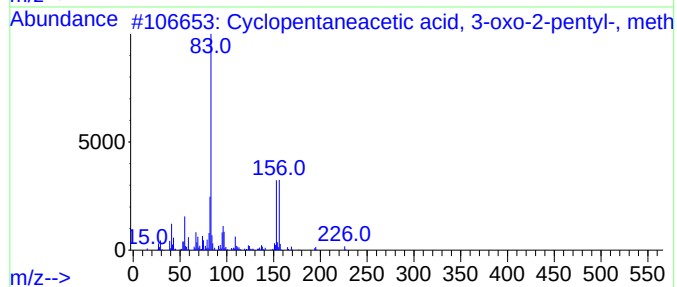
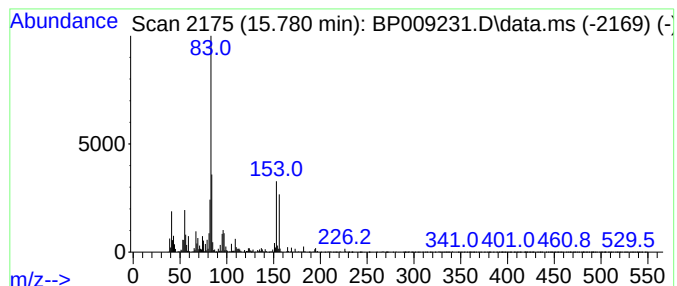
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP020722.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Cyclopentaneacetic acid, 3-... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD		R.T.	
15.780	4.91 ng	297791	Acenaphthene-d10		14.410	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Cyclopentaneacetic acid, 3-oxo-2...		226	C13H22O3	024851-98-7	89
2	Isocitronellol		156	C10H20O	018479-52-2	47
3	1,7-Octadiene, 2,3,3-trimethyl-		152	C11H20	1000150-47-7	38
4	Oxalic acid, cyclohexyl dodecyl ...		340	C20H36O4	1000309-31-4	38
5	But-2-enoylamide, 3-methyl-N-(2-...		329	C22H35NO	1000451-25-7	37



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP022122\
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Acq On : 21 Feb 2022 13:13
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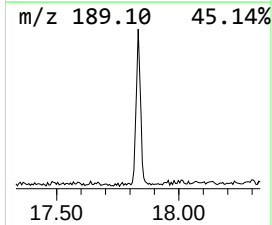
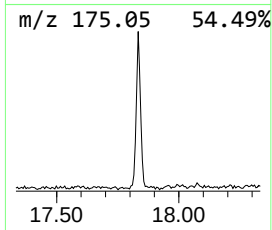
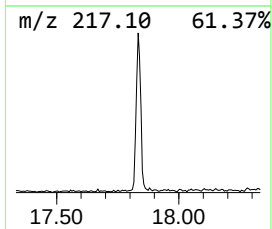
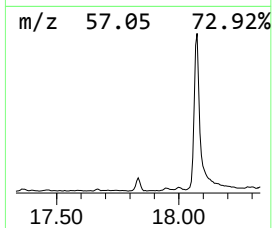
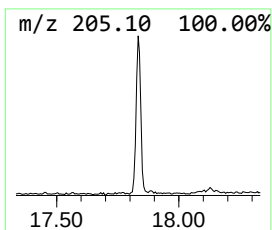
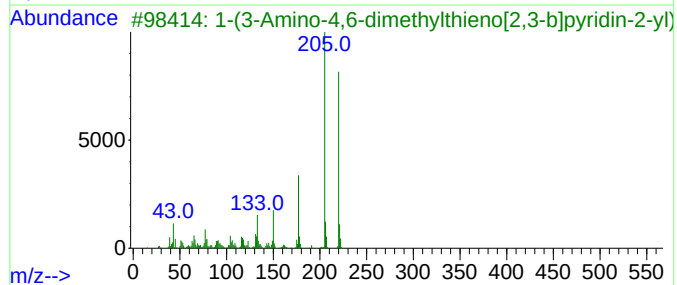
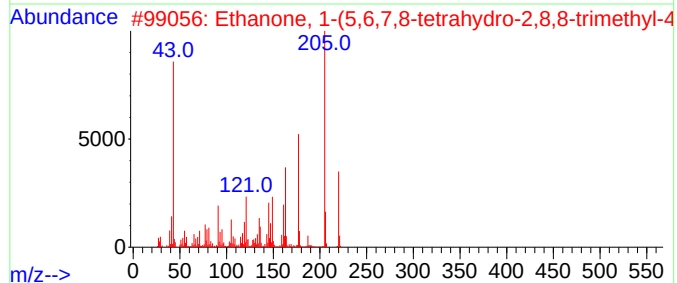
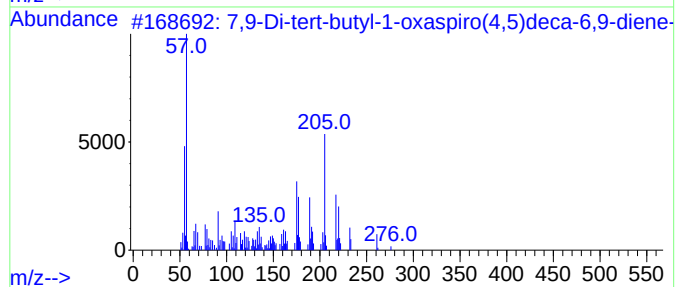
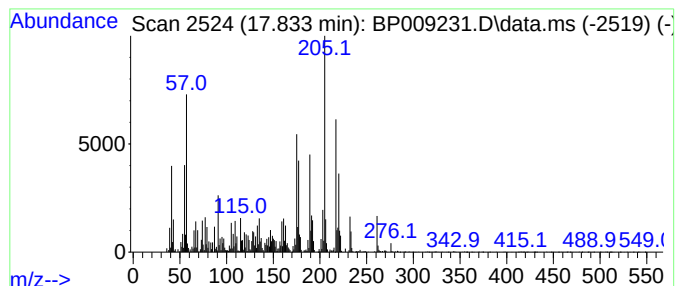
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP020722.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 7,9-Di-tert-butyl-1-oxaspiro... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.	
17.833	2.80 ng	220882	Phenanthrene-d10	17.174	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	7,9-Di-tert-butyl-1-oxaspiro(4,5...	276	C17H24O3	082304-66-3	99
2	Ethanone, 1-(5,6,7,8-tetrahydro-...	220	C14H20O2	071596-88-8	64
3	1-(3-Amino-4,6-dimethylthieno[2,...	220	C11H12N2OS	052505-42-7	50
4	Butylated Hydroxytoluene	220	C15H24O	000128-37-0	42
5	2,5-Cyclohexadien-1-one, 2,6-bis...	232	C16H24O	006738-27-8	42



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP022122\
Data File : BP009231.D
Acq On : 21 Feb 2022 13:13
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Misc :
ALS Vial : 4 Sample Multiplier: 1

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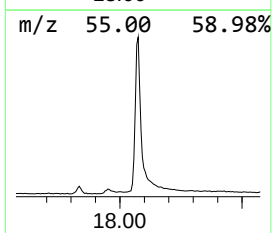
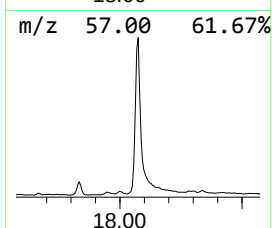
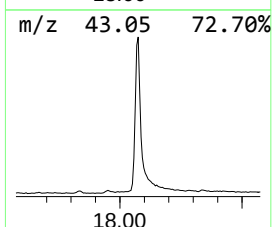
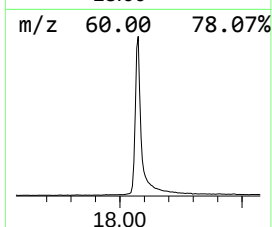
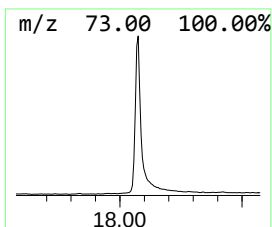
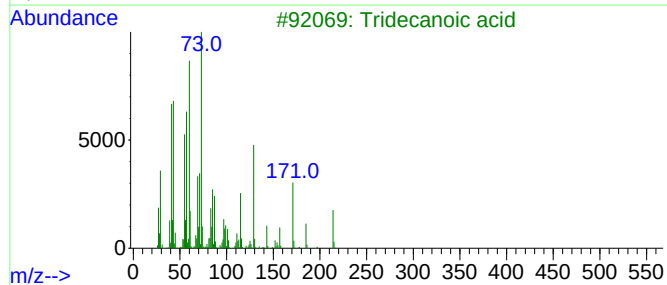
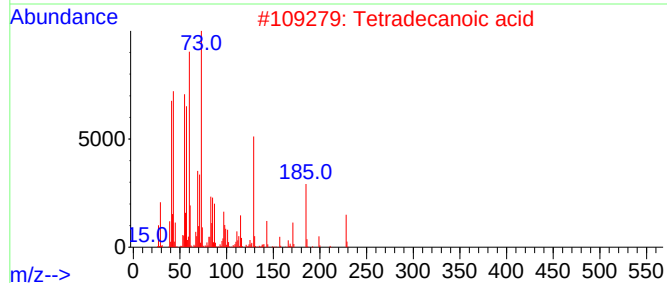
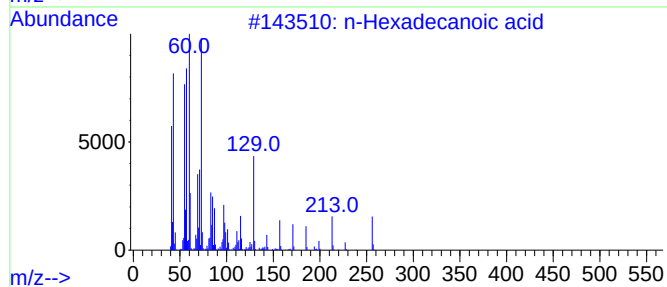
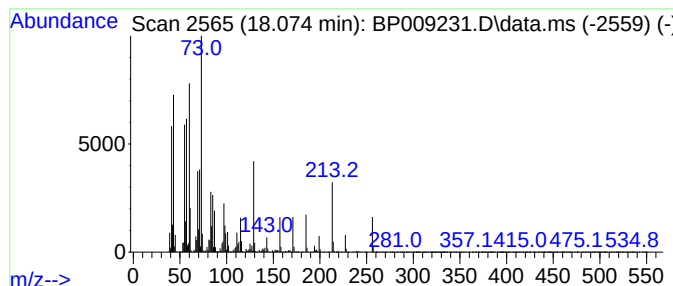
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 n-Hexadecanoic acid Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.074	43.42 ng	3424840	Phenanthrene-d10	17.174

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2		Tetradecanoic acid	228	C14H28O2	000544-63-8	89
3		Tridecanoic acid	214	C13H26O2	000638-53-9	86
4		Undecanoic acid	186	C11H22O2	000112-37-8	83
5		Pentadecanoic acid	242	C15H30O2	001002-84-2	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP022122\
Data File : BP009231.D
Acq On : 21 Feb 2022 13:13
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Sample : N1599-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
SPN-4

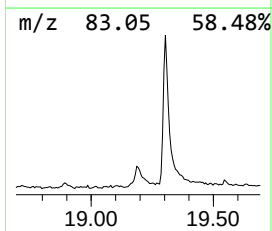
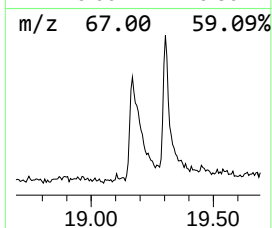
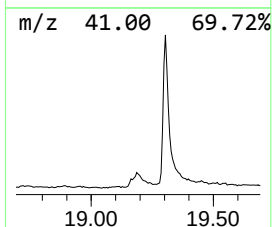
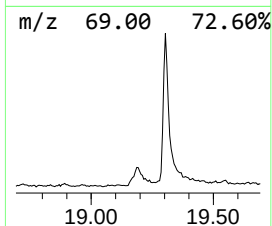
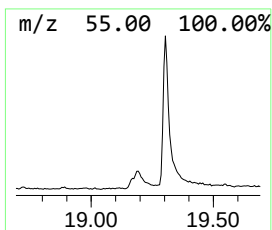
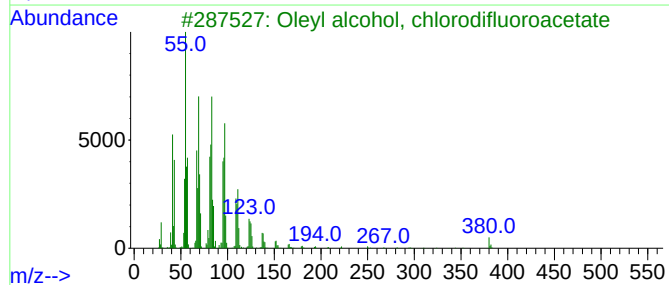
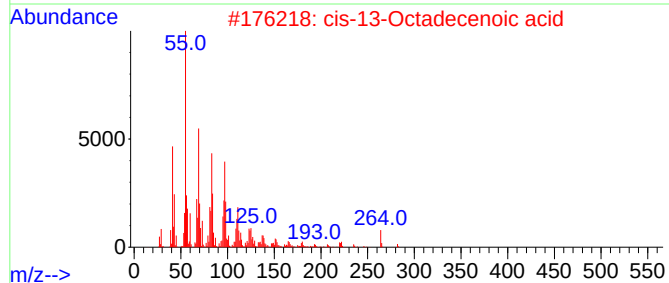
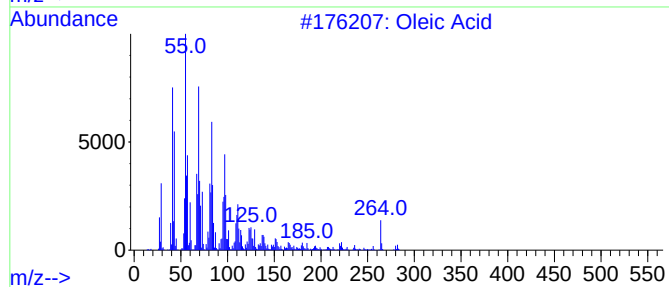
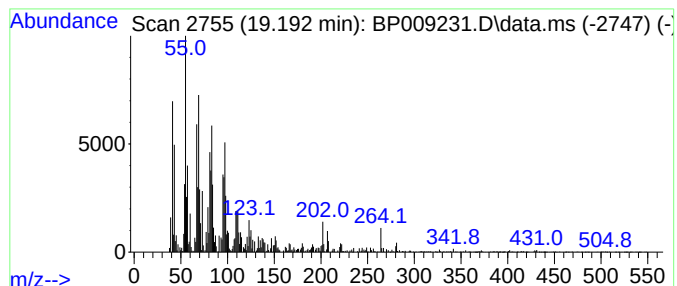
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Oleic Acid Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.192	5.07 ng	399515	Phenanthrene-d10	17.174

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Oleic Acid	282	C18H34O2	000112-80-1	91
2			cis-13-Octadecenoic acid	282	C18H34O2	013126-39-1	90
3			Oleyl alcohol, chlorodifluoroace...	380	C20H35ClF2O2	1000447-47-7	90
4			n-Propyl 11-octadecenoate	324	C21H40O2	1000336-71-7	87
5			trans-13-Octadecenoic acid	282	C18H34O2	000693-71-0	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP022122\
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Acq On : 21 Feb 2022 13:13
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Misc :
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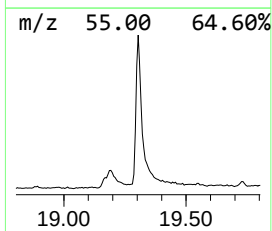
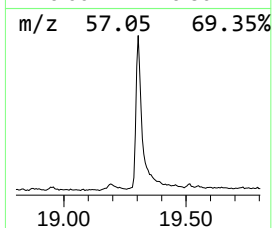
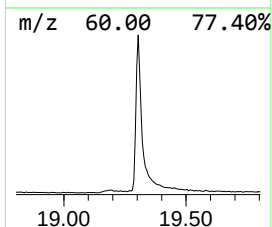
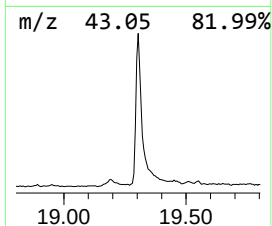
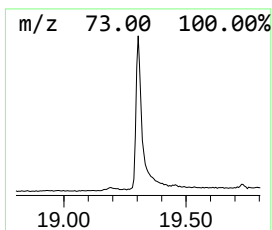
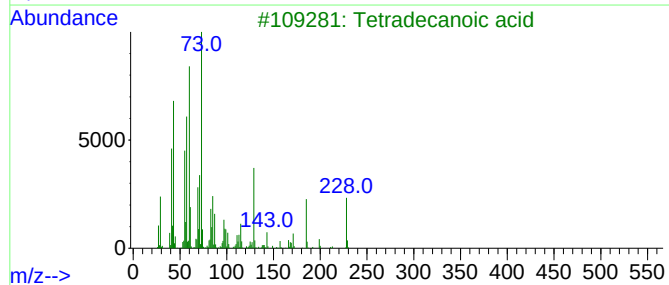
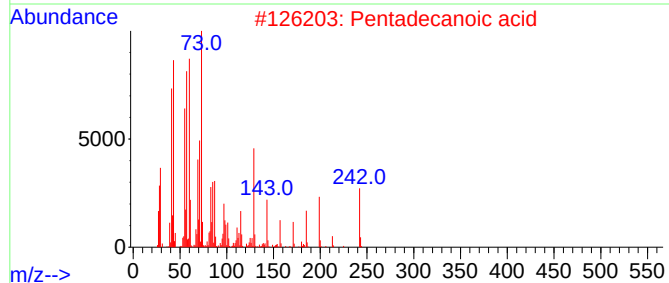
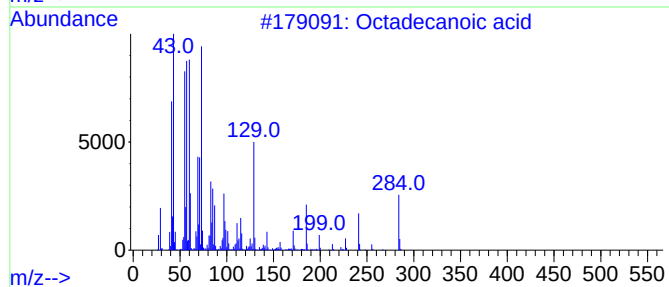
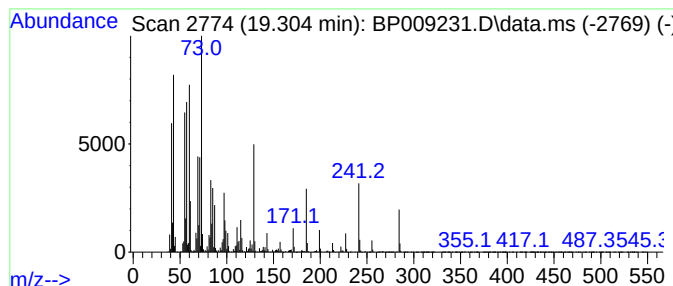
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Octadecanoic acid Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.304	20.32 ng	2281740	Chrysene-d12	21.268

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	86
3			Tetradecanoic acid	228	C14H28O2	000544-63-8	74
4			Octadecanoic acid, 2-(2-hydroxye...	372	C22H44O4	000106-11-6	72
5			Tridecanoic acid	214	C13H26O2	000638-53-9	70



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP022122\
Data File : BP009231.D
Acq On : 21 Feb 2022 13:13
Operator : CG/JU
Sample : N1599-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
SPN-4

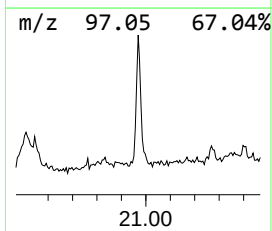
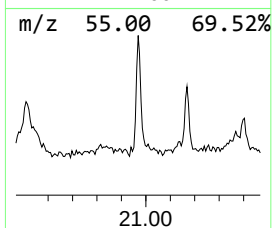
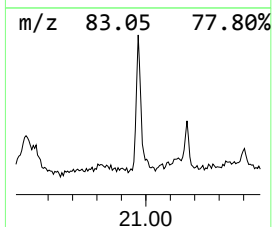
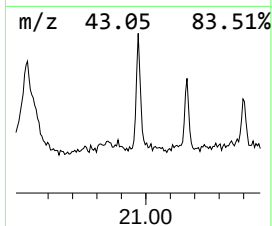
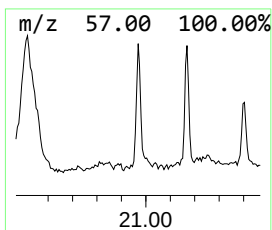
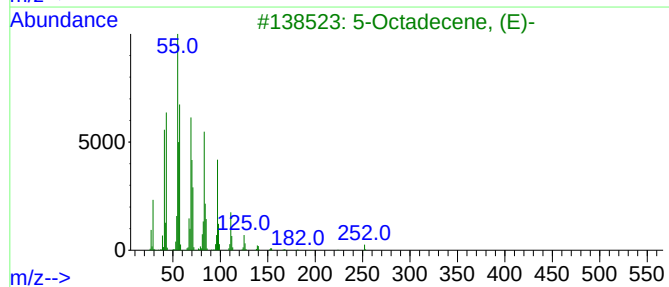
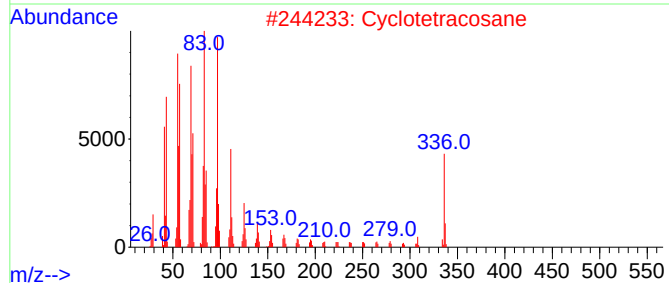
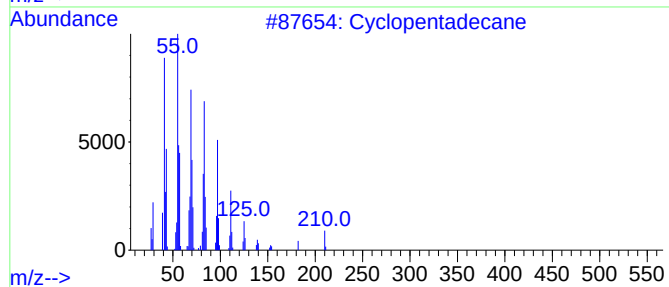
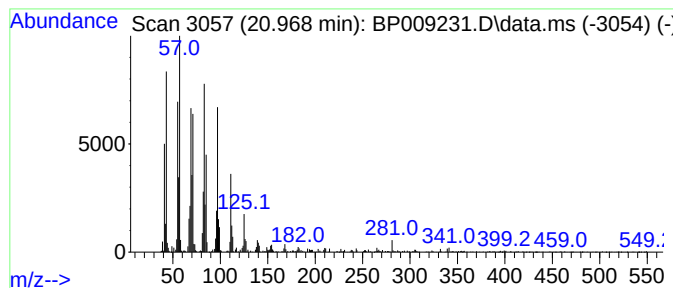
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Cyclopentadecane Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.968	2.02 ng	226736	Chrysene-d12	21.268

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cyclopentadecane	210	C15H30	000295-48-7	98
2		Cyclotetracosane	336	C24H48	000297-03-0	97
3		5-Octadecene, (E)-	252	C18H36	007206-21-5	93
4		Pentadecafluorooctanoic acid, oc...	666	C26H37F15O2	1000406-04-8	93
5		Eicosyl heptafluorobutyrate	494	C24H41F7O2	1000351-83-5	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP022122\
Data File : BP009231.D
Acq On : 21 Feb 2022 13:13
Operator : CG/JU
Sample : N1599-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
SPN-4

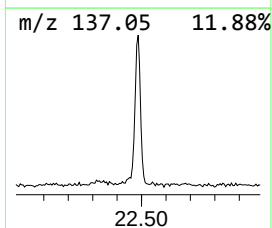
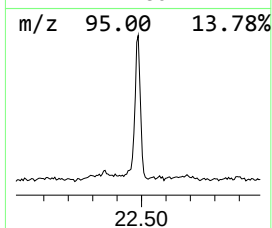
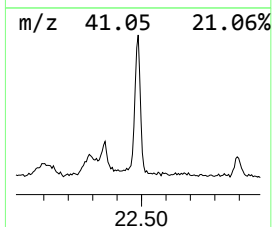
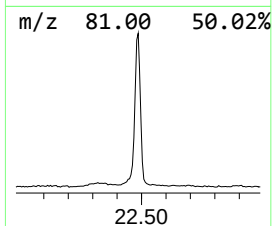
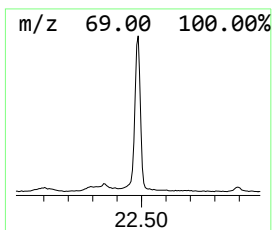
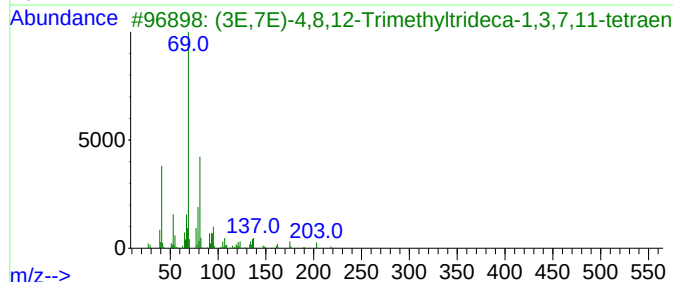
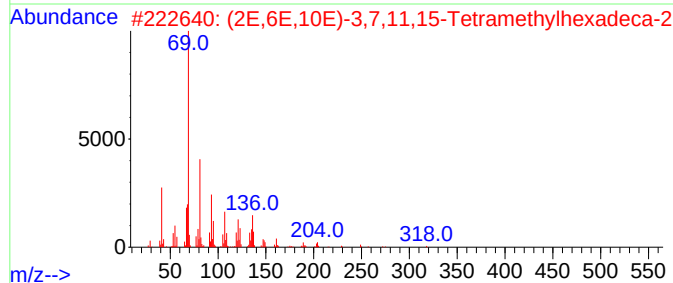
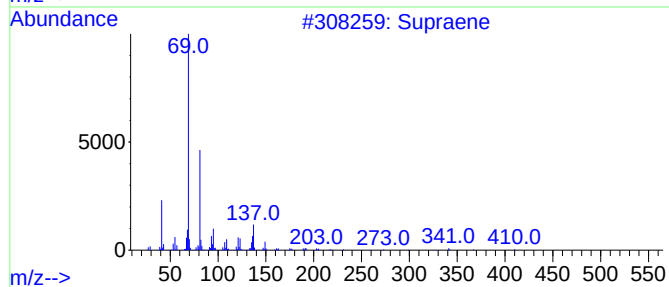
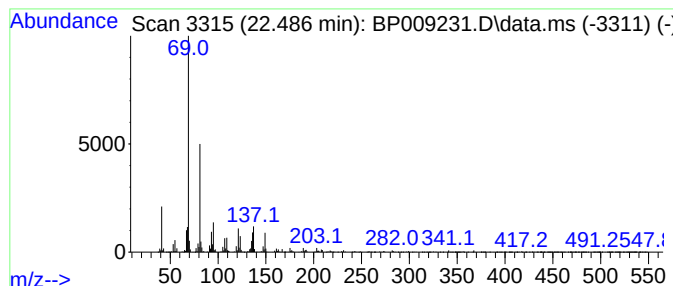
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Supraene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.486	5.69 ng	839491	Perylene-d12	23.651

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Supraene	410	C30H50	007683-64-9	91
2			(2E,6E,10E)-3,7,11,15-Tetramethy...	318	C21H34O2	125456-63-5	72
3			(3E,7E)-4,8,12-Trimethyltrideca...	218	C16H26	062235-06-7	72
4			Trifluoroacetyl-lavandulol	250	C12H17F3O2	028673-24-7	64
5			4,8,12-Tetradecatrienal, 5,9,13-...	248	C17H28O	066408-55-7	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP022122\
 Data File : BP009231.D
 Acq On : 21 Feb 2022 13:13
 Operator : CG/JU
 Sample : N1599-01
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SPN-4

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP020722.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
Cyclopentanone	4.246	10.9	ng	291875	1	7.763	537774	20.0
2-Propanol, 1-(...	7.316	5.6	ng	150189	1	7.763	537774	20.0
Ethyl 2,5,8,11,...	7.369	5.6	ng	149699	1	7.763	537774	20.0
2-Propanol, 1-(...	7.557	10.0	ng	268052	1	7.763	537774	20.0
1-Hexanol, 2-et...	7.828	10.1	ng	271629	1	7.763	537774	20.0
2,2,4-Trimethyl...	12.763	3.8	ng	227612	3	14.410	1211940	20.0
Phenol, 4-(1,1-...	13.263	18.5	ng	1122860	3	14.410	1211940	20.0
Cyclopentaneace...	15.780	4.9	ng	297791	3	14.410	1211940	20.0
7,9-Di-tert-but...	17.833	2.8	ng	220882	4	17.174	1577460	20.0
n-Hexadecanoic ...	18.074	43.4	ng	3424840	4	17.174	1577460	20.0
Oleic Acid	19.192	5.1	ng	399515	4	17.174	1577460	20.0
Octadecanoic acid	19.304	20.3	ng	2281740	5	21.268	2245260	20.0
Cyclopentadecane	20.968	2.0	ng	226736	5	21.268	2245260	20.0
Supraene	22.486	5.7	ng	839491	6	23.651	2948970	20.0