

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP021722\
 Data File : BP009195.D
 Acq On : 17 Feb 2022 15:02
 Operator : CG/JU
 Sample : N1575-01
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 COP-1R

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: BP009195.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.975	164	168	175	rVB	140461	197781	1.72%	0.332%
2	4.263	212	217	226	rVB	234594	341598	2.97%	0.573%
3	4.905	320	326	338	rBV2	109694	196872	1.71%	0.330%
4	5.375	400	406	412	rBV	2190888	3674983	31.92%	6.161%
5	5.787	470	476	481	rBV	151562	237413	2.06%	0.398%
6	6.975	672	678	698	rVB	1282240	2740978	23.81%	4.595%
7	7.328	733	738	743	rBV	154297	246092	2.14%	0.413%
8	7.381	743	747	751	rVV	161557	275672	2.39%	0.462%
9	7.422	751	754	768	rVB	83113	159296	1.38%	0.267%
10	7.575	773	780	794	rBV	217663	447820	3.89%	0.751%
11	7.775	806	814	821	rBV	761558	1306446	11.35%	2.190%
12	7.846	821	826	830	rVV	164767	294663	2.56%	0.494%
13	7.887	830	833	847	rVB	79326	146141	1.27%	0.245%
14	8.940	1005	1012	1036	rBV	2284538	4232334	36.76%	7.095%
15	10.575	1282	1290	1297	rBV	871216	1706053	14.82%	2.860%
16	12.775	1654	1664	1678	rBV	138432	270228	2.35%	0.453%
17	13.039	1700	1709	1734	rBV	7202408	11394342	98.96%	19.101%
18	13.275	1740	1749	1766	rVB	722149	1232318	10.70%	2.066%
19	14.316	1918	1926	1937	rBV2	67536	216700	1.88%	0.363%
20	14.422	1937	1944	1951	rVV	1532127	2382117	20.69%	3.993%
21	15.245	2077	2084	2094	rVB2	133161	241980	2.10%	0.406%
22	15.792	2171	2177	2183	rBV2	110078	176939	1.54%	0.297%
23	15.922	2192	2199	2217	rBV	4869869	7791408	67.67%	13.061%
24	17.180	2407	2413	2428	rBV	1536962	2528982	21.96%	4.239%
25	17.845	2521	2526	2531	rBV2	117719	166181	1.44%	0.279%
26	18.080	2561	2566	2575	rBV	267088	511692	4.44%	0.858%
27	19.316	2772	2776	2788	rBV	161814	334439	2.90%	0.561%
28	19.745	2843	2849	2861	rBV	9081804	11514164	100.00%	19.302%
29	21.057	3068	3072	3080	rBV	262878	437241	3.80%	0.733%
30	21.280	3106	3110	3122	rBV	1430756	2061448	17.90%	3.456%
31	23.668	3510	3516	3533	rVB2	875097	2189403	19.01%	3.670%

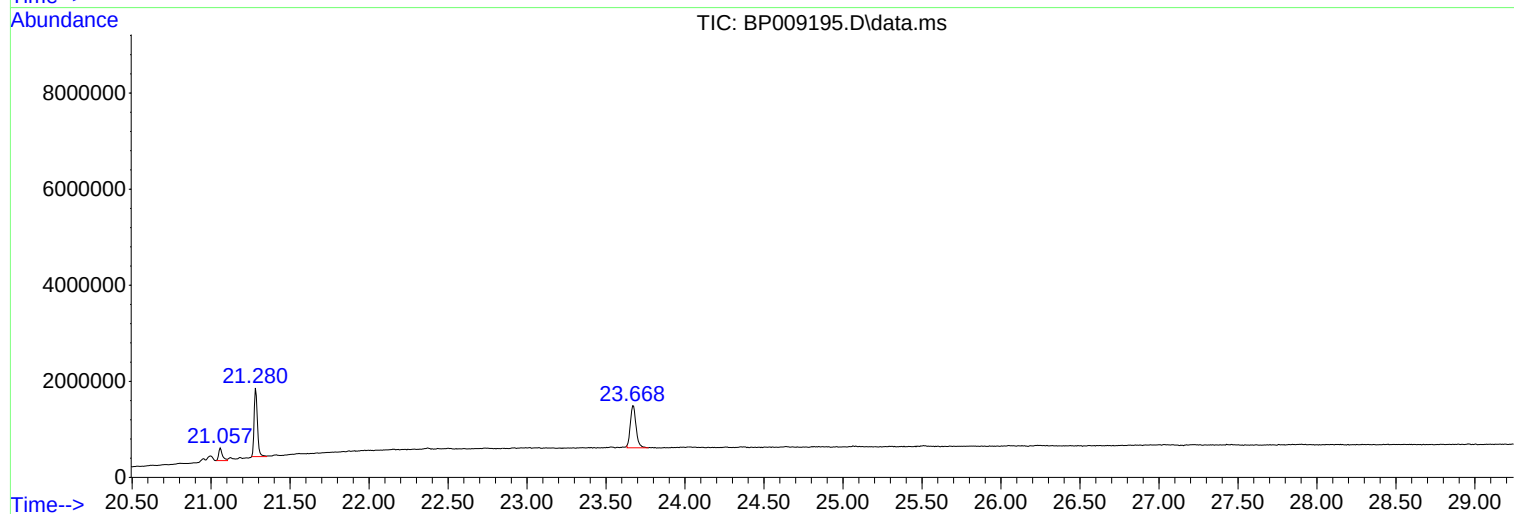
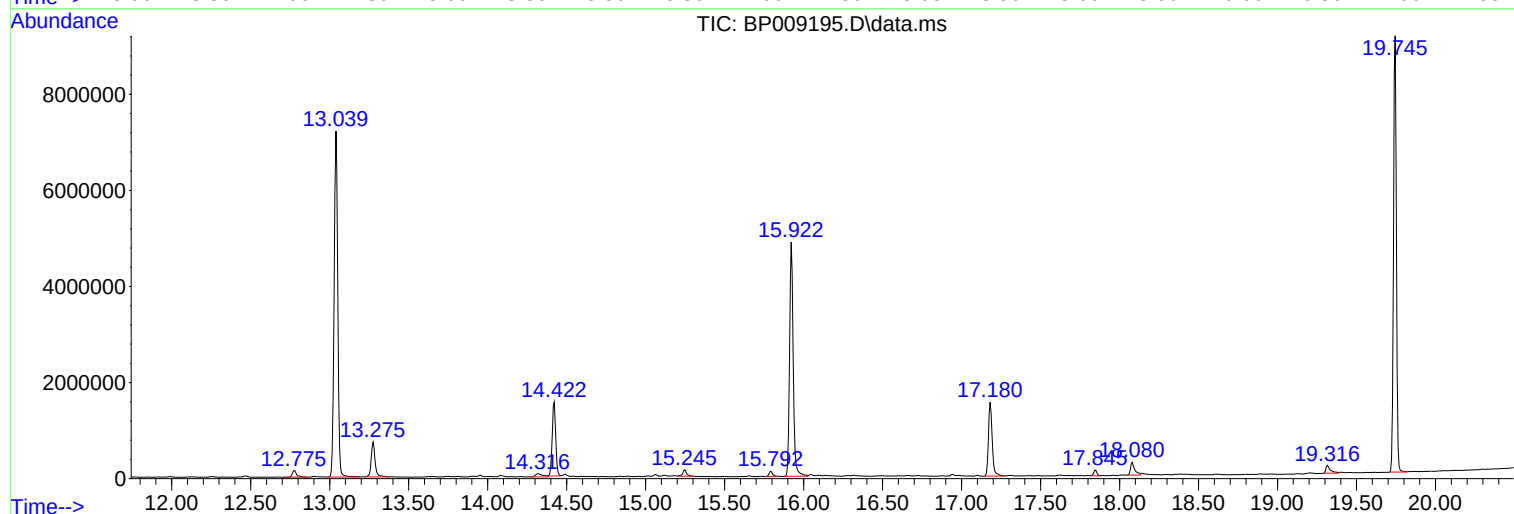
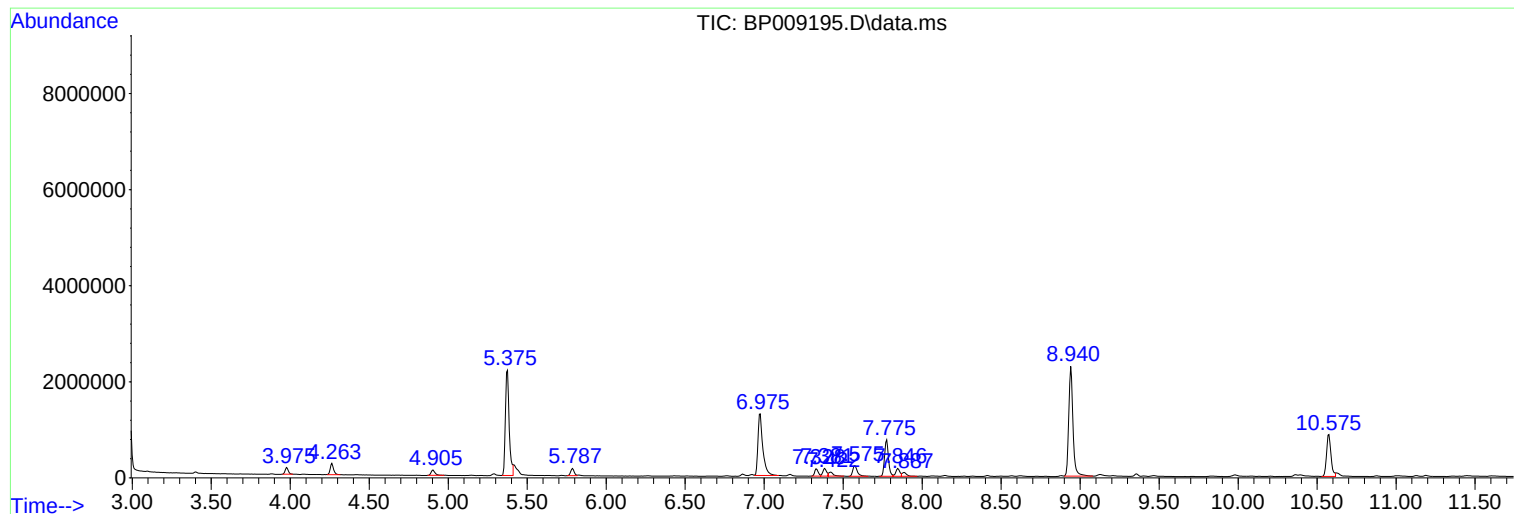
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP020722.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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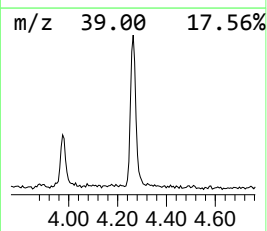
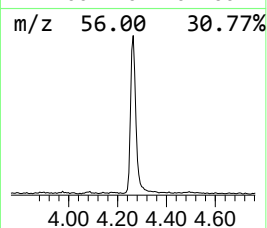
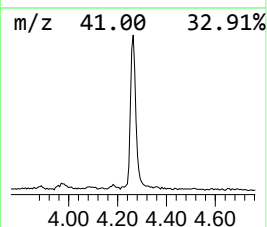
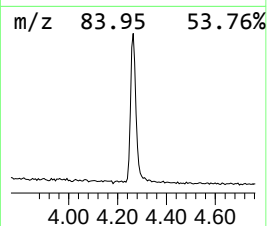
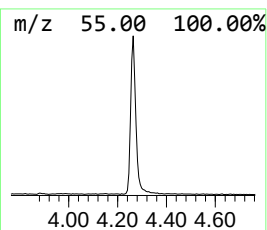
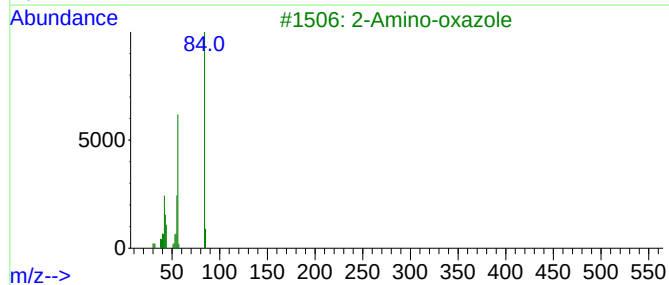
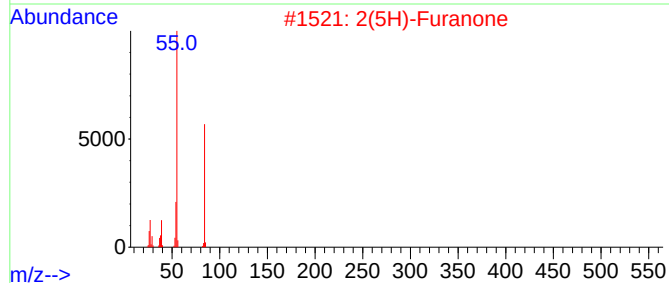
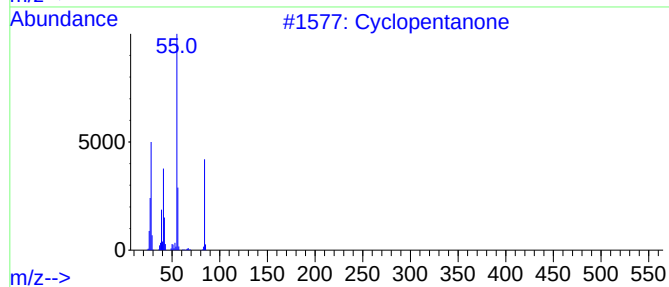
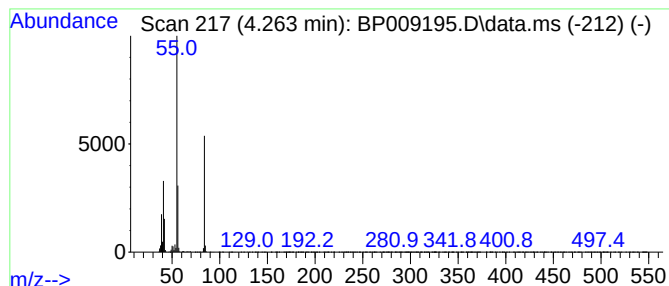
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP020722.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 Cyclopentanone Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.263	5.23 ng	341598	1,4-Dichlorobenzene-d4	7.775

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



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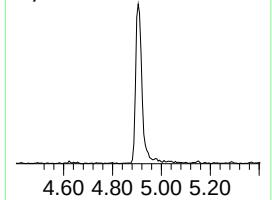
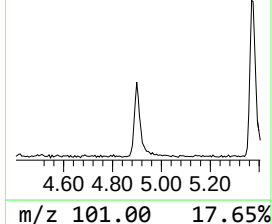
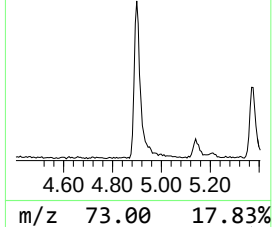
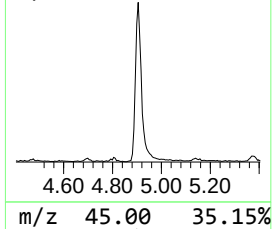
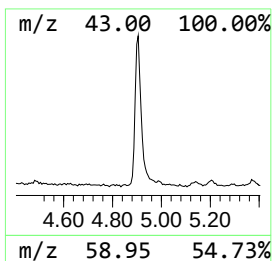
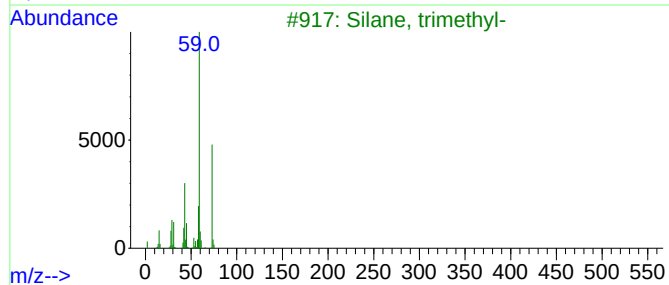
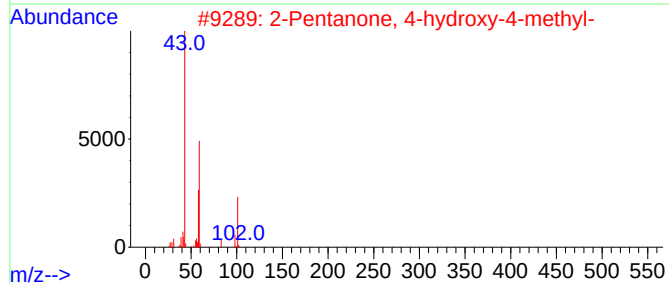
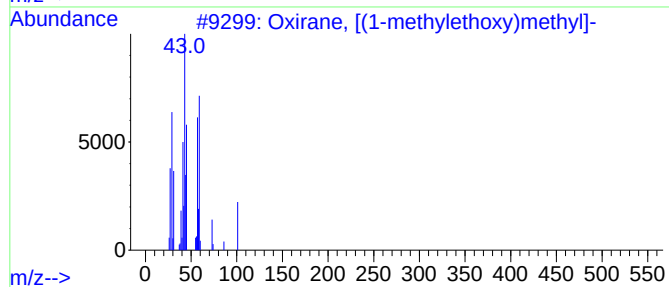
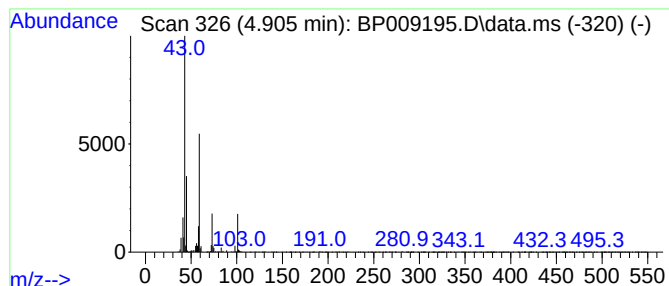
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP020722.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 3 Oxirane, [(1-methylethoxy)m... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.905	3.01 ng	196872	1,4-Dichlorobenzene-d4	7.775

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Oxirane, [(1-methylethoxy)methyl]-	116	C6H12O2	004016-14-2	50
2			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	47
3			Silane, trimethyl-	74	C3H10Si	000993-07-7	43
4			Ethane, 1-ethoxy-1-methoxy-	104	C5H12O2	010471-14-4	43
5			Tripropylene glycol, propyl ethe...	348	C18H40O4Si	1000367-07-2	38



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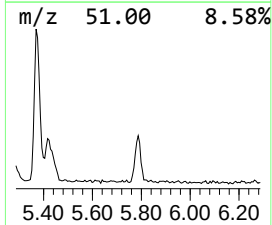
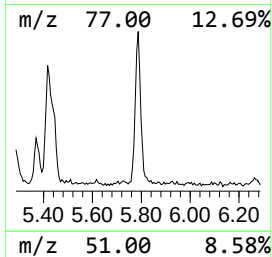
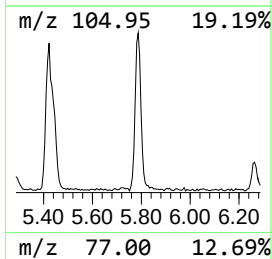
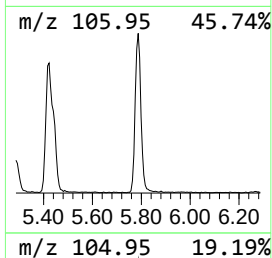
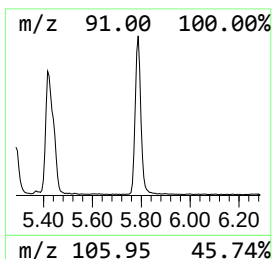
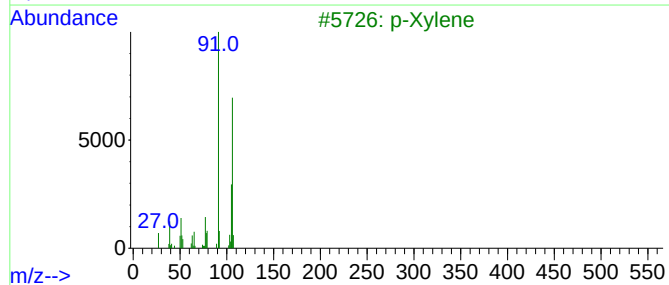
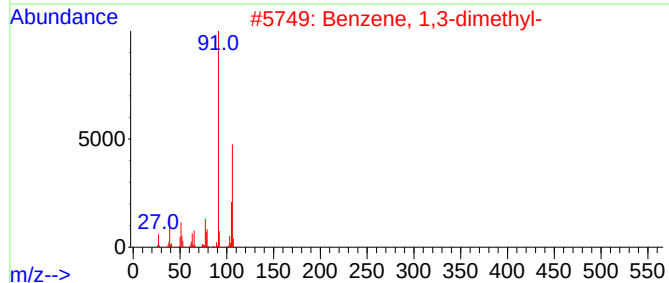
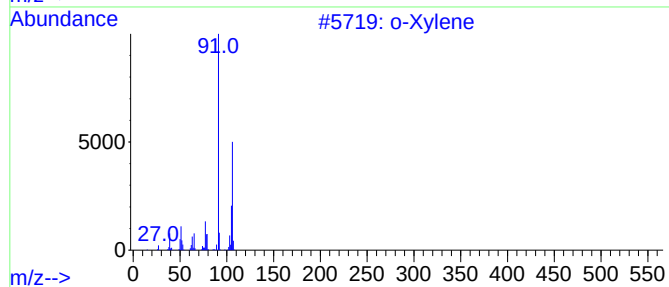
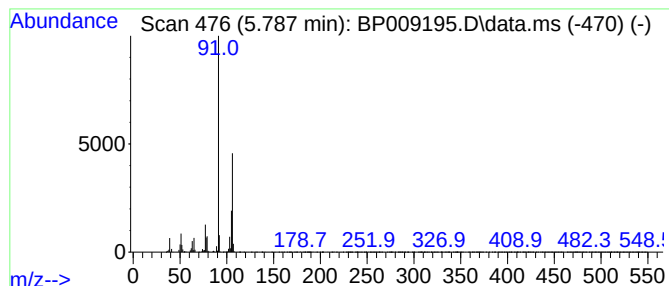
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 4 o-Xylene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.787	3.63 ng	237413	1,4-Dichlorobenzene-d4	7.775

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			o-Xylene	106	C8H10	000095-47-6	97
2			Benzene, 1,3-dimethyl-	106	C8H10	000108-38-3	97
3			p-Xylene	106	C8H10	000106-42-3	95
4			1,3-Cyclopentadiene, 5-(1-methyl...	106	C8H10	002175-91-9	87
5			Ethylbenzene	106	C8H10	000100-41-4	83



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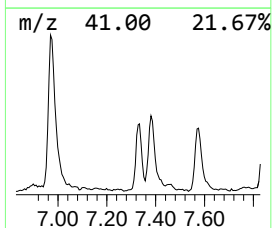
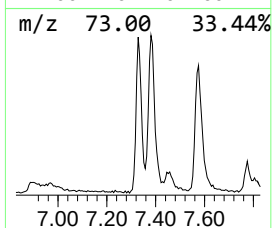
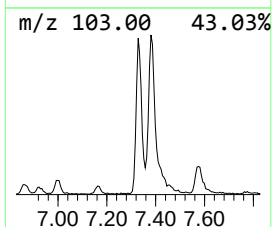
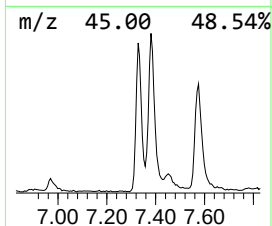
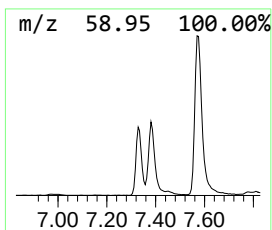
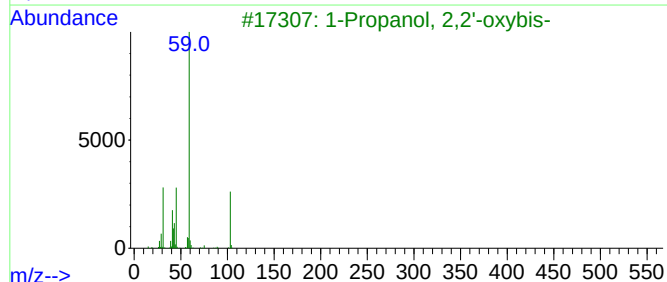
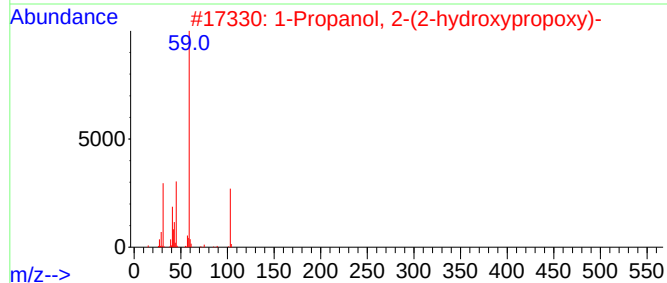
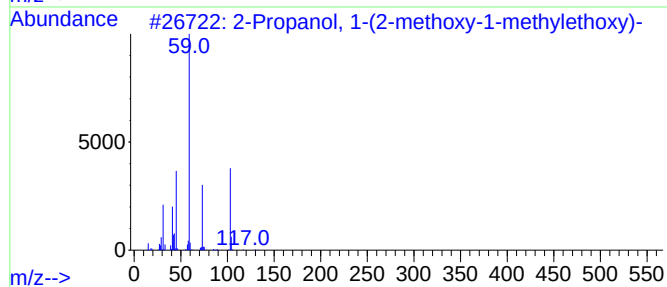
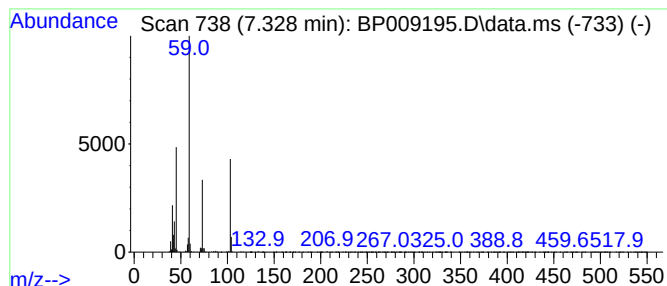
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 2-Propanol, 1-(2-methoxy-1-... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.328	3.77 ng	246092	1,4-Dichlorobenzene-d4	7.775

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-hydroxypropoxy)-	134	C6H14O3	000106-62-7	59
3			1-Propanol, 2,2'-oxybis-	134	C6H14O3	000108-61-2	59
4			Dipropylene glycol (isomer 2)	134	C6H14O3	1000506-25-4	53
5			Dipropylene glycol (isomer 3)	134	C6H14O3	1000506-25-5	53



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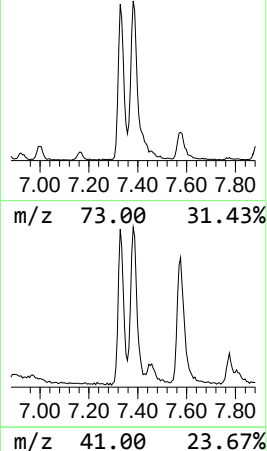
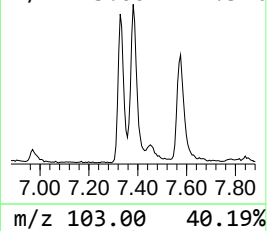
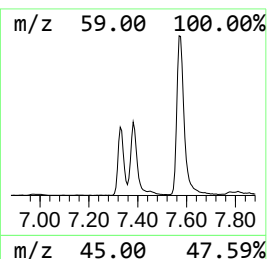
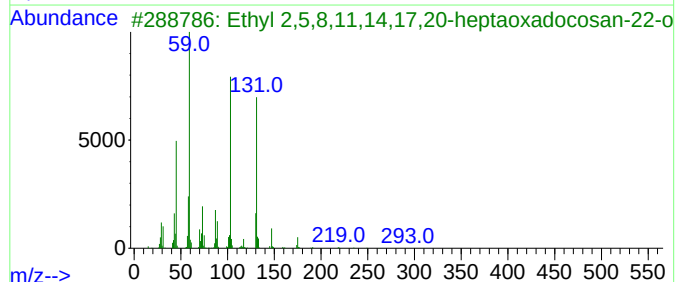
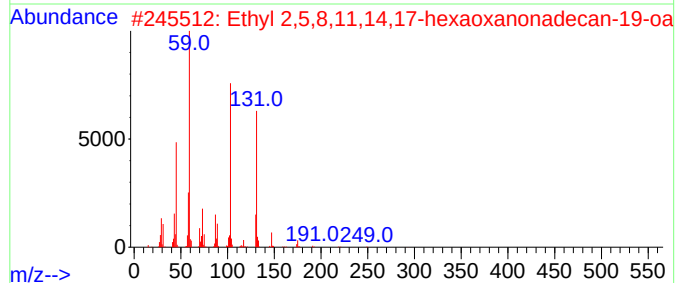
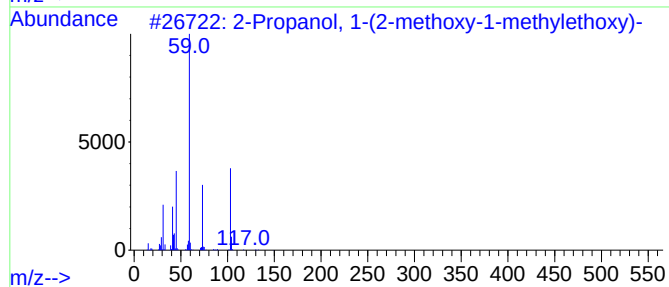
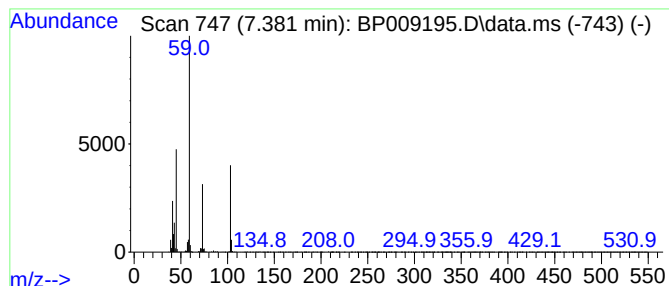
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TIC Library : C:\Database\NIST20.L
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Peak Number 6 Ethyl 2,5,8,11,14,17-hexaox... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.381	4.22 ng	275672	1,4-Dichlorobenzene-d4	7.775

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Propanol, 1-(2-methoxy-1-methy...	148	C7H16O3	020324-32-7	90		
2	Ethyl 2,5,8,11,14,17-hexaoxanona...	338	C15H30O8	1000366-80-6	72		
3	Ethyl 2,5,8,11,14,17,20-heptaoxa...	382	C17H34O9	1000366-80-5	56		
4	Dipropylene glycol (isomer 2)	134	C6H14O3	1000506-25-4	53		
5	Dipropylene glycol (isomer 3)	134	C6H14O3	1000506-25-5	53		



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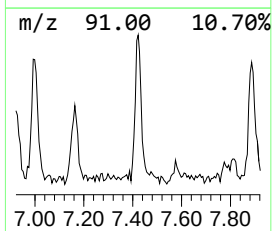
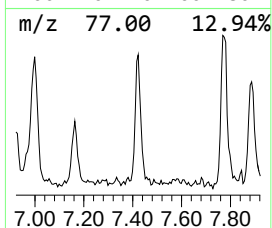
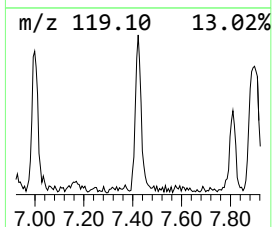
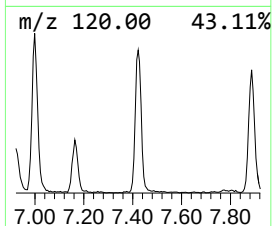
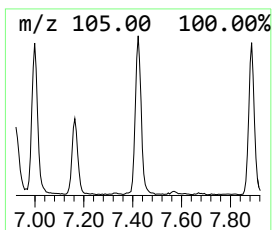
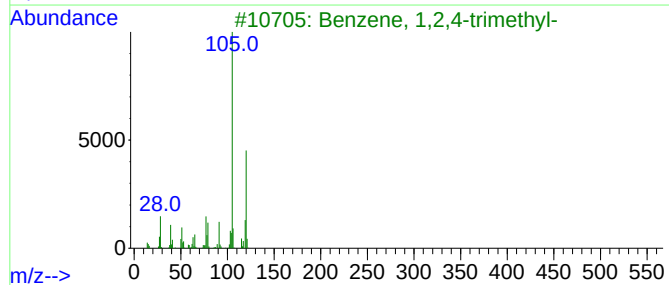
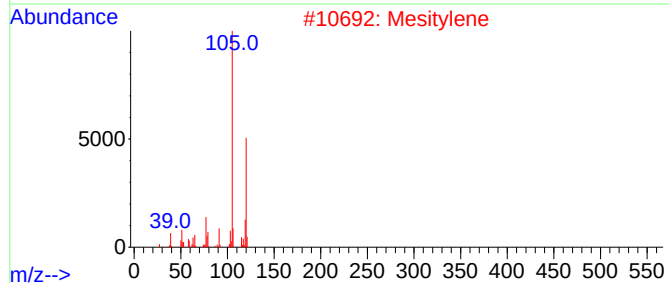
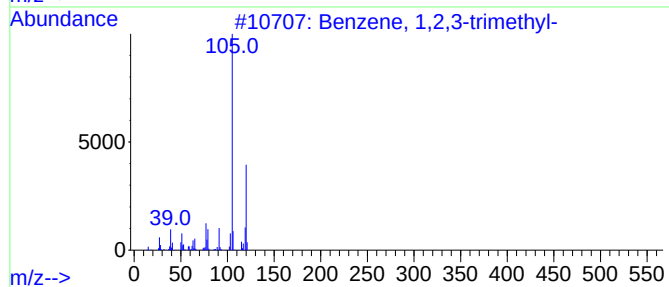
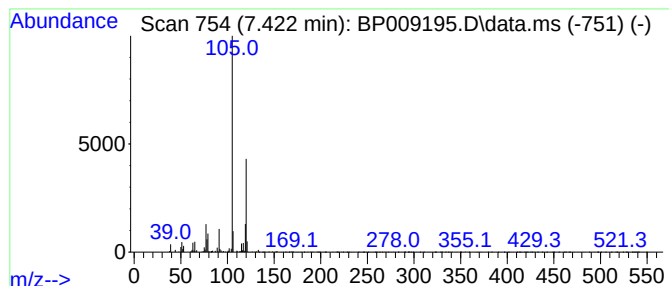
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Benzene, 1,2,3-trimethyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.422	2.44 ng	159296	1,4-Dichlorobenzene-d4	7.775

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	94
2			Mesitylene	120	C9H12	000108-67-8	93
3			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	90
4			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	90
5			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP021722\
Data File : BP009195.D
Acq On : 17 Feb 2022 15:02
Operator : CG/JU
Sample : N1575-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

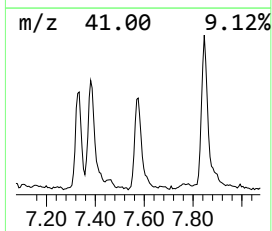
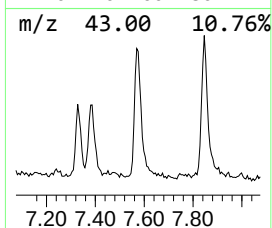
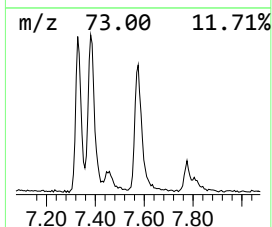
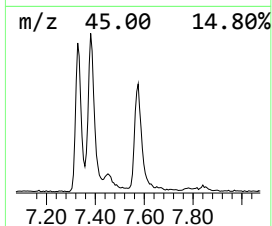
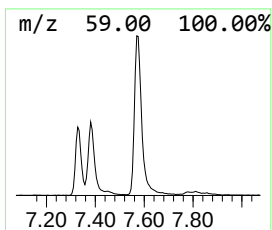
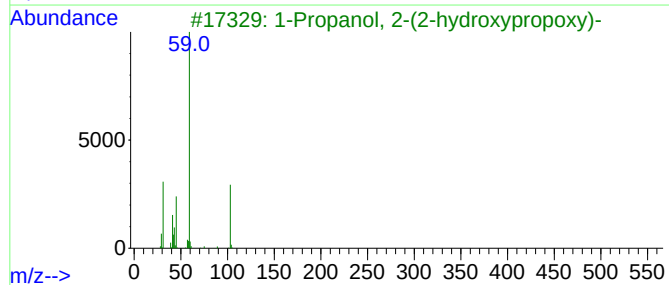
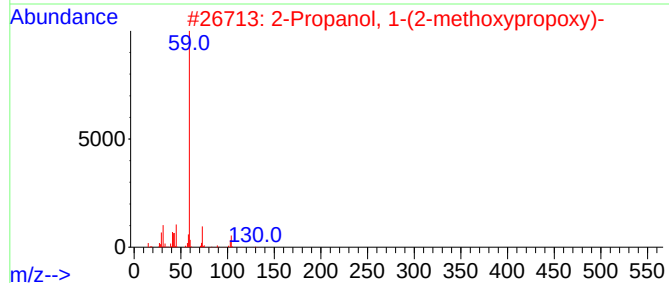
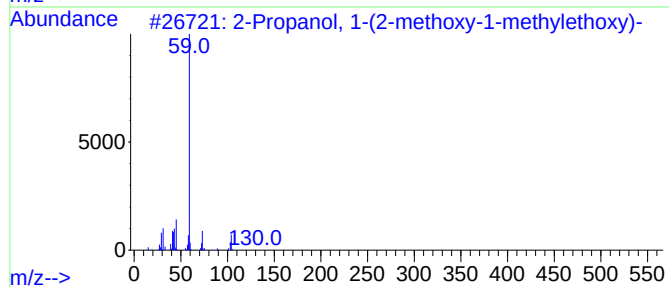
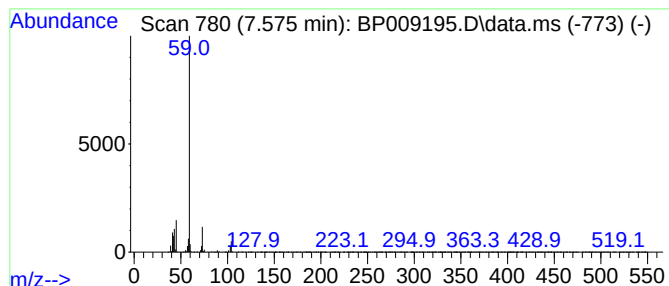
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP020722.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD			R.T.
7.575	6.86 ng	447820	1,4-Dichlorobenzene-d4			7.775
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	1-Propanol,	2-(2-hydroxypropoxy)-	134	C6H14O3	000106-62-7	64
4	2-Propanol,	1-[1-methyl-2-(2-pro...	174	C9H18O3	055956-25-7	43
5	2-Propanol,	1-(2-butoxy-1-methyl...	190	C10H22O3	029911-28-2	40



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP021722\
Data File : BP009195.D
Acq On : 17 Feb 2022 15:02
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Sample : N1575-01
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ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
COP-1R

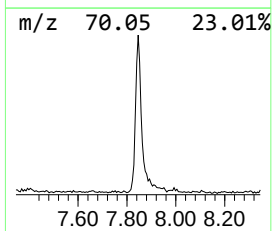
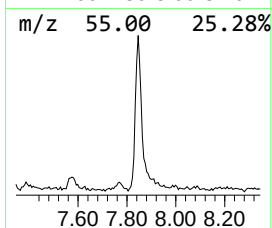
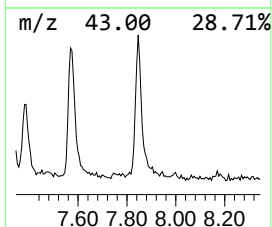
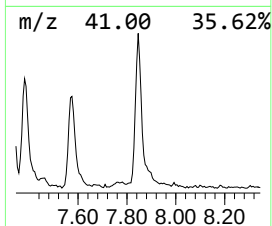
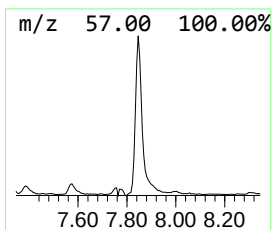
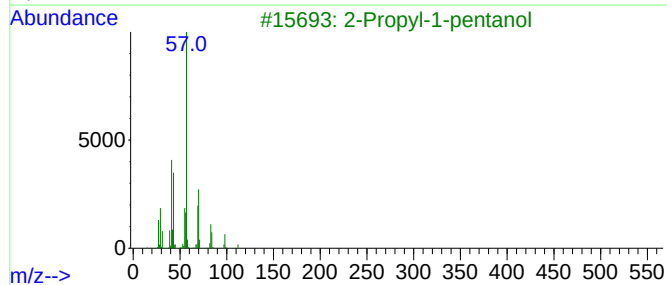
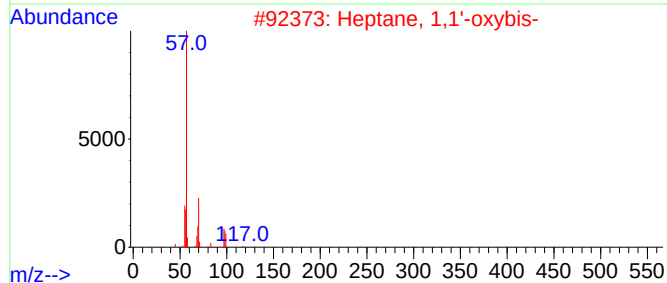
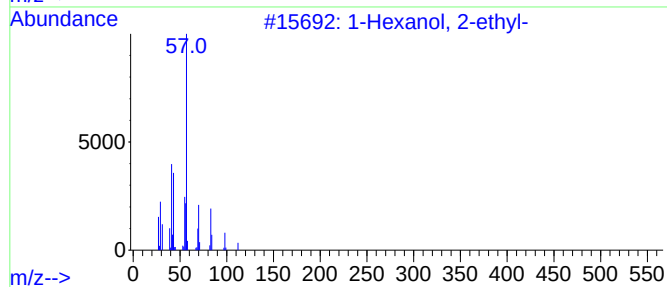
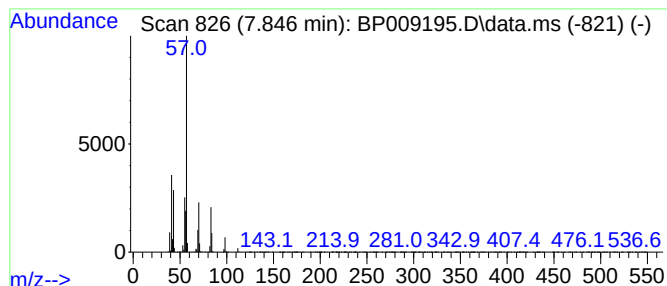
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 1-Hexanol, 2-ethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.846	4.51 ng	294663	1,4-Dichlorobenzene-d4	7.775

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	83
2			Heptane, 1,1'-oxybis-	214	C14H30O	000629-64-1	53
3			2-Propyl-1-pentanol	130	C8H18O	058175-57-8	53
4			Chloroacetic acid, 2-ethylhexyl ...	206	C10H19ClO2	005345-58-4	45
5			Carbonic acid, isobutyl 2-ethylh...	230	C13H26O3	1000383-13-3	42



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP021722\
Data File : BP009195.D
Acq On : 17 Feb 2022 15:02
Operator : CG/JU
Sample : N1575-01
Misc :
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Instrument :
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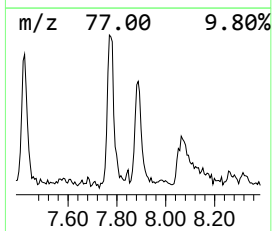
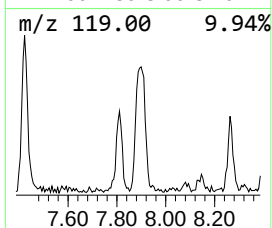
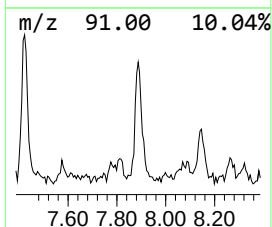
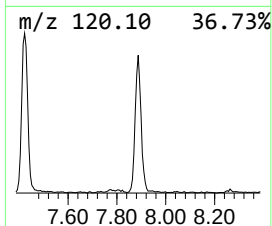
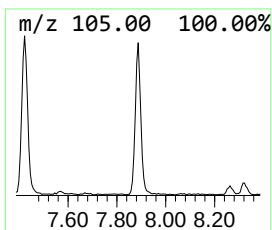
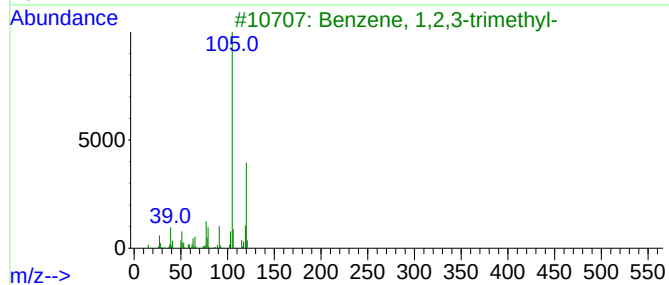
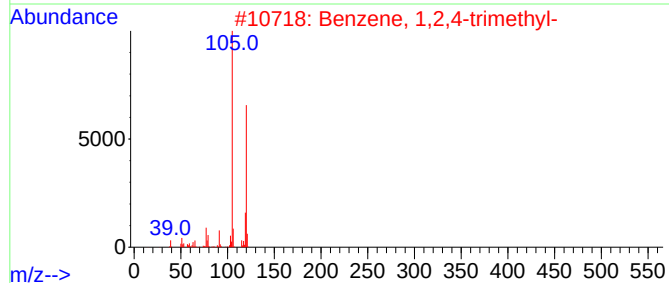
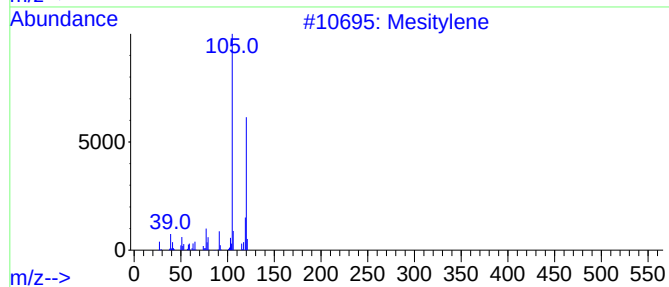
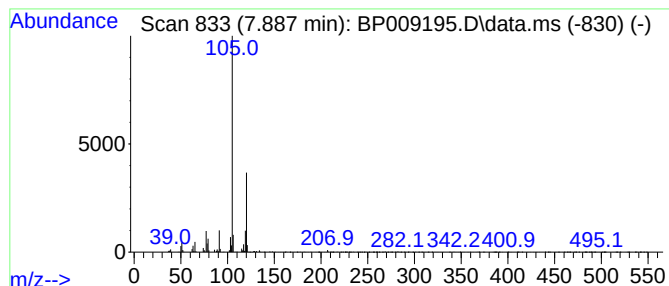
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Mesitylene Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.887	2.24 ng	146141	1,4-Dichlorobenzene-d4	7.775

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Mesitylene	120	C9H12	000108-67-8	93
2			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
3			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
4			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	90
5			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP021722\
Data File : BP009195.D
Acq On : 17 Feb 2022 15:02
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Sample : N1575-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

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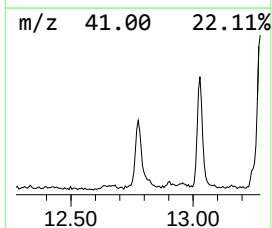
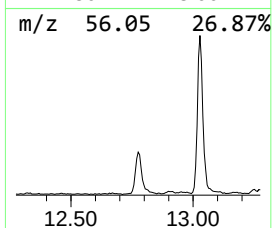
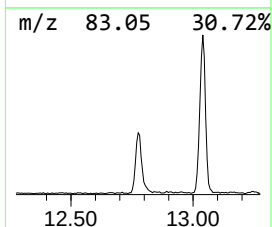
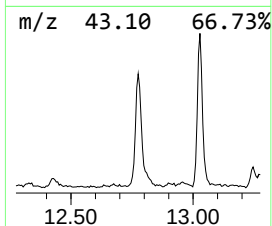
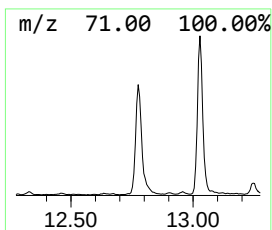
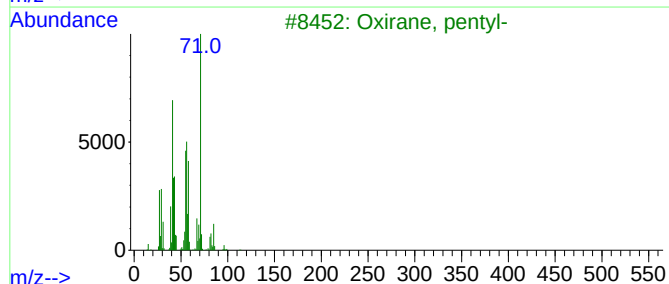
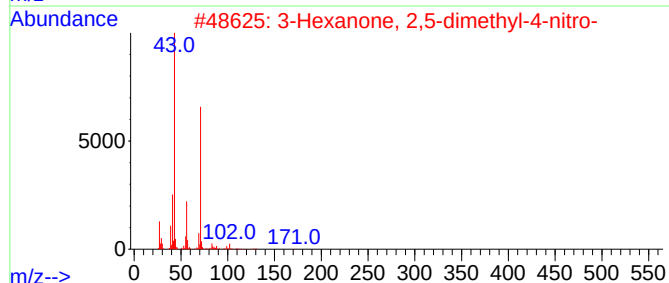
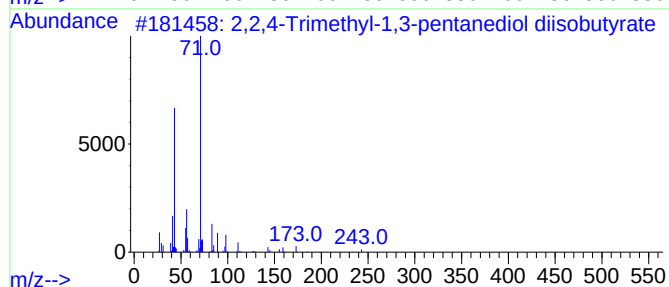
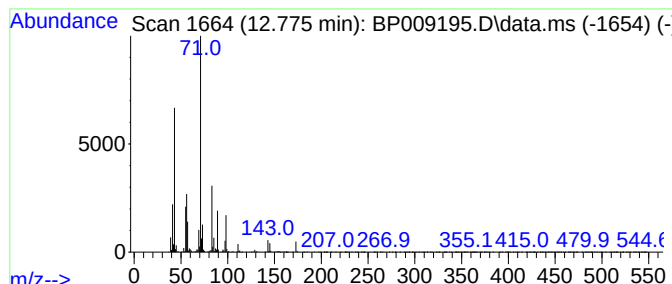
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 2,2,4-Trimethyl-1,3-pentane... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.775	2.27 ng	270228	Acenaphthene-d10	14.422

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2,2,4-Trimethyl-1,3-pentanediol ...	286	C16H30O4	006846-50-0	72
2			3-Hexanone, 2,5-dimethyl-4-nitro-	173	C8H15NO3	059906-54-6	43
3			Oxirane, pentyl-	114	C7H14O	005063-65-0	43
4			Butyric acid, 3-tridecyl ester	270	C17H34O2	1000280-56-8	43
5			2-METHYLHEPT-1-EN-3-OL	128	C8H16O	013019-19-7	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP021722\
Data File : BP009195.D
Acq On : 17 Feb 2022 15:02
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Sample : N1575-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

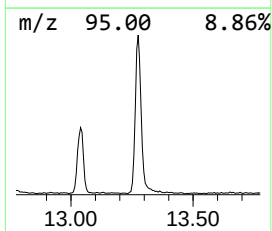
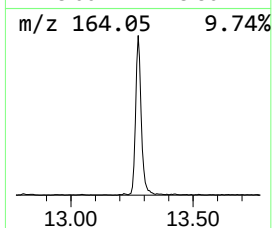
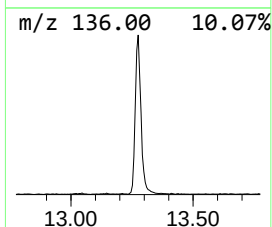
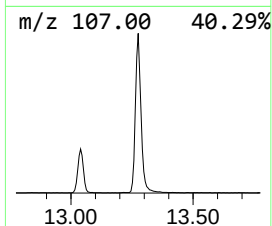
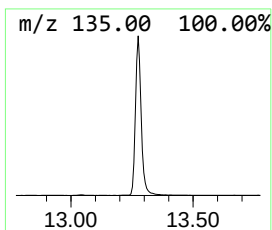
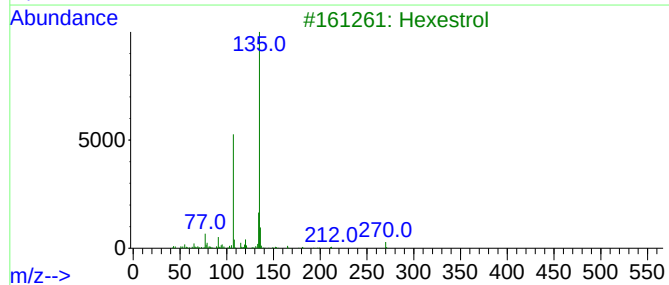
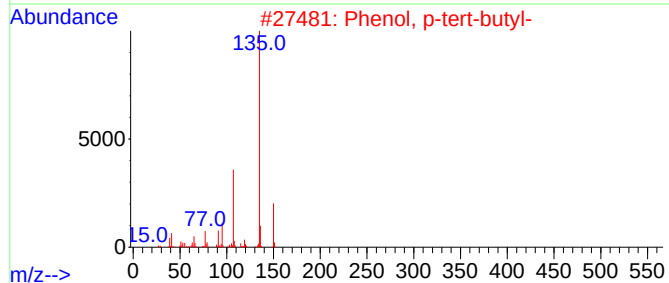
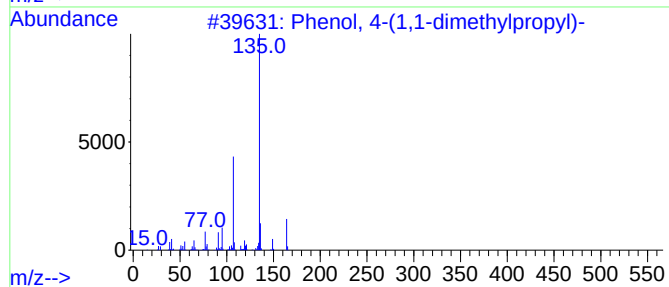
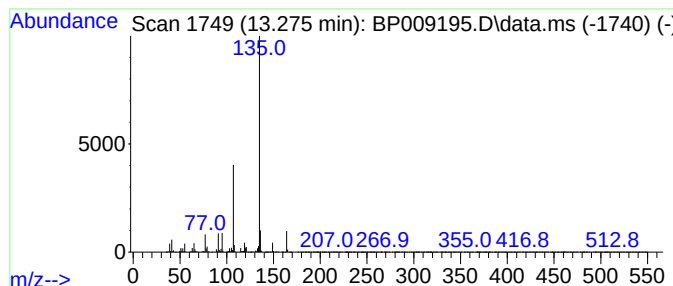
Instrument :
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ClientSampleId :
COP-1R

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 12 Phenol, 4-(1,1-dimethylprop... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD			R.T.
13.275	10.35 ng	1232320	Acenaphthene-d10			14.422
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Phenol, 4-(1,1-dimethylpropyl)-		164	C11H16O	000080-46-6	95
2	Phenol, p-tert-butyl-		150	C10H14O	000098-54-4	80
3	Hexestrol		270	C18H22O2	000084-16-2	72
4	Hexestrol, O-trifluoroacetyl-		366	C20H21F3O3	1000365-31-7	72
5	2-tert-Butylphenol, O-isopropyllo...		236	C14H20O3	1000487-38-4	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP021722\
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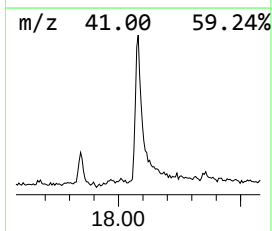
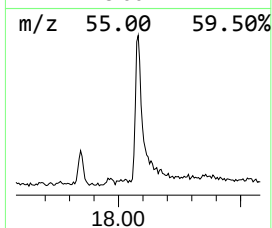
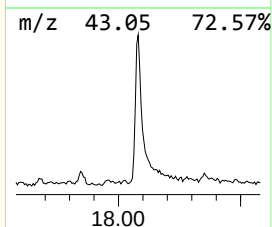
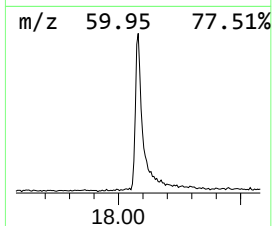
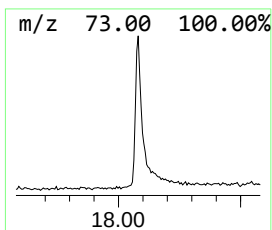
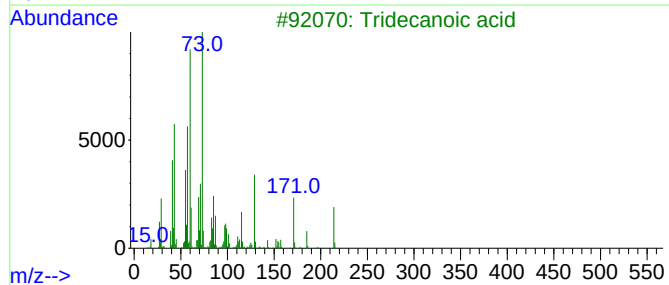
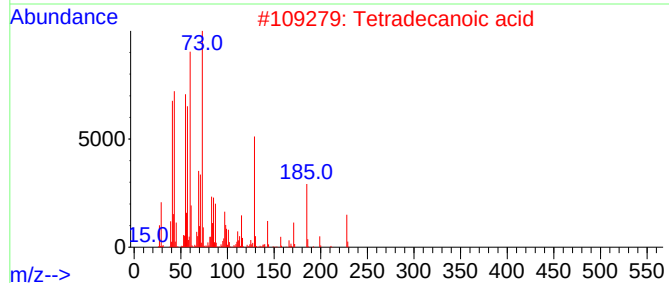
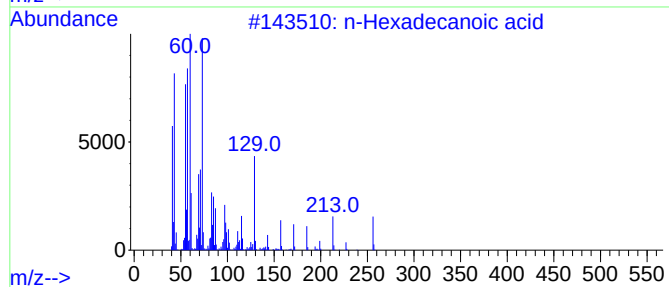
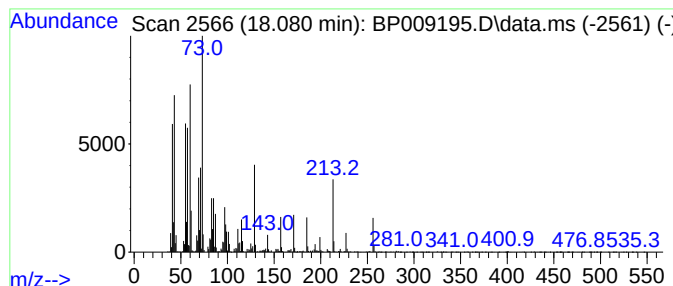
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 n-Hexadecanoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD		R.T.	
18.080	4.05 ng	511692	Phenanthrene-d10		17.180	
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	93
3	Tridecanoic acid		214	C13H26O2	000638-53-9	89
4	Pentadecanoic acid		242	C15H30O2	001002-84-2	87
5	Octadecanoic acid		284	C18H36O2	000057-11-4	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP021722\
Data File : BP009195.D
Acq On : 17 Feb 2022 15:02
Operator : CG/JU
Sample : N1575-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
COP-1R

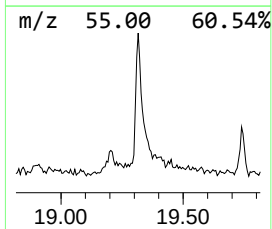
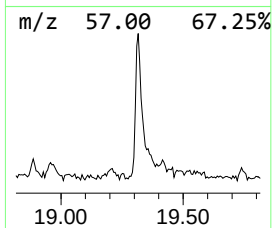
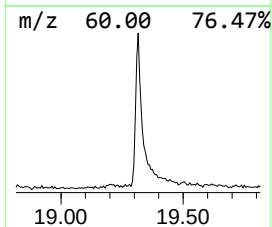
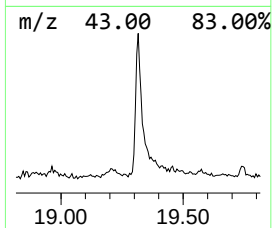
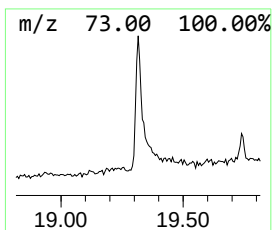
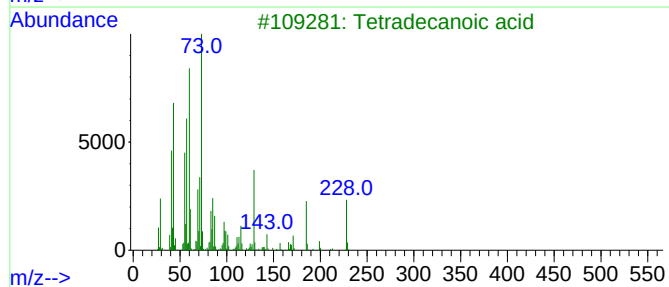
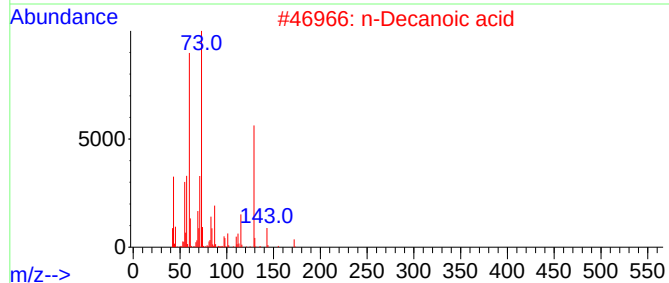
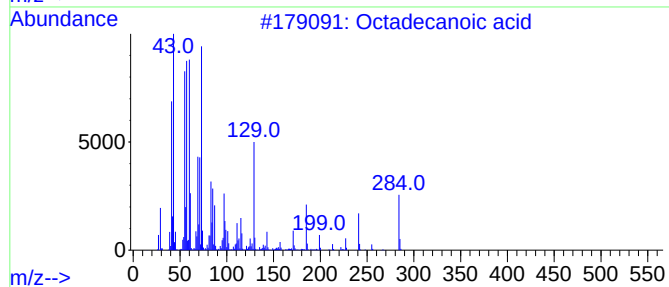
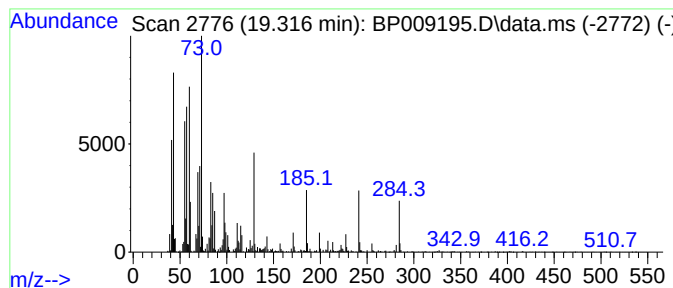
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Octadecanoic acid Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.316	3.24 ng	334439	Chrysene-d12	21.280

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid	284	C18H36O2	000057-11-4	99
2			n-Decanoic acid	172	C10H20O2	000334-48-5	78
3			Tetradecanoic acid	228	C14H28O2	000544-63-8	74
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	64
5			Pentadecanoic acid	242	C15H30O2	001002-84-2	58



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP021722\
Data File : BP009195.D
Acq On : 17 Feb 2022 15:02
Operator : CG/JU
Sample : N1575-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
COP-1R

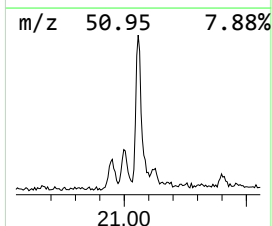
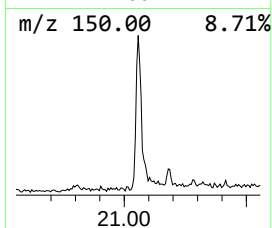
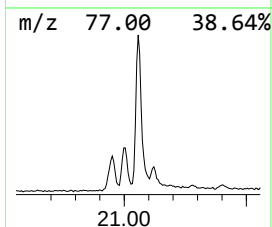
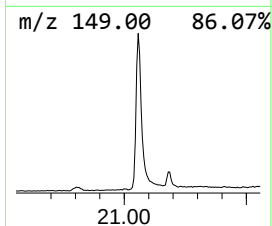
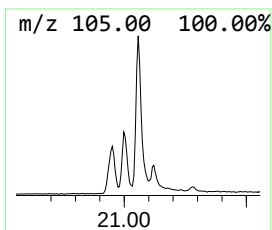
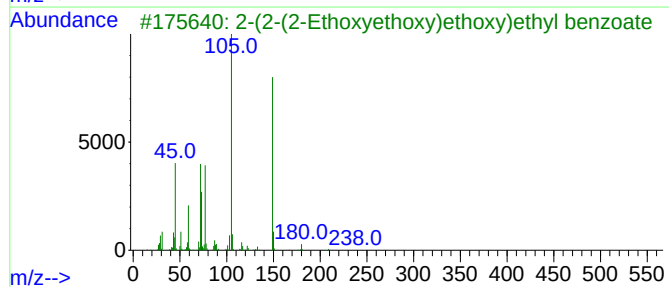
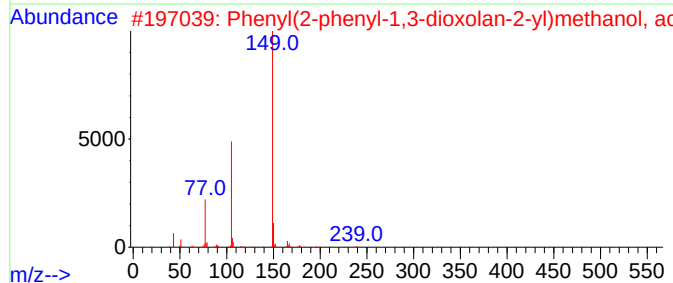
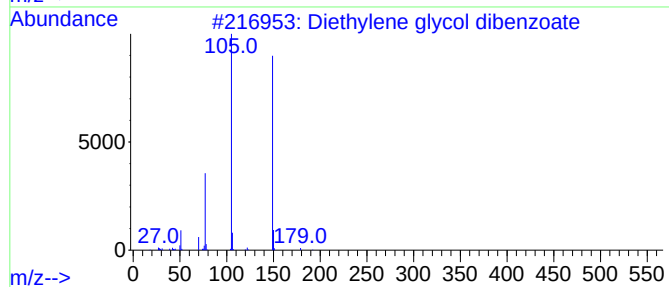
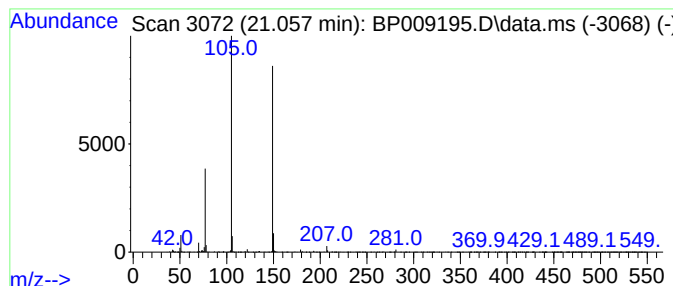
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP020722.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Diethylene glycol dibenzoate Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.057	4.24 ng	437241	Chrysene-d12	21.280

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Diethylene glycol dibenzoate	314	C18H18O5	000120-55-8	90
2			Phenyl(2-phenyl-1,3-dioxolan-2-yl)methanol, ac	298	C18H18O4	006963-16-2	64
3			2-(2-(2-Ethoxyethoxy)ethoxy)ethyl benzoate	282	C15H22O5	1000366-98-8	64
4			2,5,8,11-Tetraoxatridecan-13-yl ...	312	C16H24O6	1000366-96-7	64
5			2-(2-(2-Methoxyethoxy)ethoxy)ethyl benzoate	268	C14H20O5	1000366-99-1	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP021722\
 Data File : BP009195.D
 Acq On : 17 Feb 2022 15:02
 Operator : CG/JU
 Sample : N1575-01
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 COP-1R

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	4.263	5.2	ng	341598	1	7.775	1306450	20.0
Oxirane, [(1-me...	4.905	3.0	ng	196872	1	7.775	1306450	20.0
o-Xylene	5.787	3.6	ng	237413	1	7.775	1306450	20.0
2-Propanol, 1-(...	7.328	3.8	ng	246092	1	7.775	1306450	20.0
Ethyl 2,5,8,11,...	7.381	4.2	ng	275672	1	7.775	1306450	20.0
Benzene, 1,2,3-...	7.422	2.4	ng	159296	1	7.775	1306450	20.0
2-Propanol, 1-(...	7.575	6.9	ng	447820	1	7.775	1306450	20.0
1-Hexanol, 2-et...	7.846	4.5	ng	294663	1	7.775	1306450	20.0
Mesitylene	7.887	2.2	ng	146141	1	7.775	1306450	20.0
2,2,4-Trimethyl...	12.775	2.3	ng	270228	3	14.422	2382120	20.0
Phenol, 4-(1,1-...	13.275	10.4	ng	1232320	3	14.422	2382120	20.0
n-Hexadecanoic ...	18.080	4.0	ng	511692	4	17.180	2528980	20.0
Octadecanoic acid	19.316	3.2	ng	334439	5	21.280	2061450	20.0
Diethylene glyc...	21.057	4.2	ng	437241	5	21.280	2061450	20.0