

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070921\
 Data File : BP006161.D
 Acq On : 09 Jul 2021 15:15
 Operator : CG/JU
 Sample : M2990-13
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 TP-J

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-BP070721.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BP006161.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.452	395	402	420	rBV	2256669	3514508	29.29%	5.372%
2	7.058	669	675	693	rBV	1983180	3471939	28.94%	5.307%
3	7.869	805	813	823	rBV	626343	1001350	8.35%	1.531%
4	9.040	1005	1012	1024	rBV	1280226	2189115	18.25%	3.346%
5	10.469	1248	1255	1265	rBV	254412	496296	4.14%	0.759%
6	10.675	1283	1290	1299	rBV	790764	1350579	11.26%	2.064%
7	13.134	1701	1708	1718	rBV	3492586	5025180	41.88%	7.681%
8	14.510	1936	1942	1951	rVB	1208125	1700274	14.17%	2.599%
9	15.869	2155	2173	2174	rBV2	376242	1437056	11.98%	2.197%
10	16.004	2190	2196	2216	rVB2	2696524	4737019	39.48%	7.241%
11	16.634	2279	2303	2324	rVB	2376994	11998023	100.00%	18.339%
12	16.951	2346	2357	2373	rVB	150826	632673	5.27%	0.967%
13	17.192	2391	2398	2405	rBV4	102818	347674	2.90%	0.531%
14	17.263	2405	2410	2418	rVB	1354321	1979011	16.49%	3.025%
15	18.151	2554	2561	2570	rBV	1447441	2398159	19.99%	3.666%
16	18.433	2605	2609	2612	rBV	130232	187989	1.57%	0.287%
17	18.475	2613	2616	2628	rVB	158611	262879	2.19%	0.402%
18	19.380	2765	2770	2788	rVB	770960	1635467	13.63%	2.500%
19	19.816	2839	2844	2852	rBV	6711830	7992662	66.62%	12.217%
20	19.922	2857	2862	2871	rVB2	264562	611606	5.10%	0.935%
21	20.316	2925	2929	2939	rVB	400549	675632	5.63%	1.033%
22	20.510	2956	2962	2972	rVB4	242060	663473	5.53%	1.014%
23	21.010	3042	3047	3051	rBV	527539	929941	7.75%	1.421%
24	21.069	3051	3057	3061	rVV	1233466	2207033	18.39%	3.373%
25	21.116	3061	3065	3073	rVB	2198217	2737127	22.81%	4.184%
26	21.345	3100	3104	3115	rBV	1695174	2176738	18.14%	3.327%
27	21.780	3174	3178	3184	rVB	405750	692103	5.77%	1.058%
28	23.745	3506	3512	3519	rVB2	1227008	2371530	19.77%	3.625%

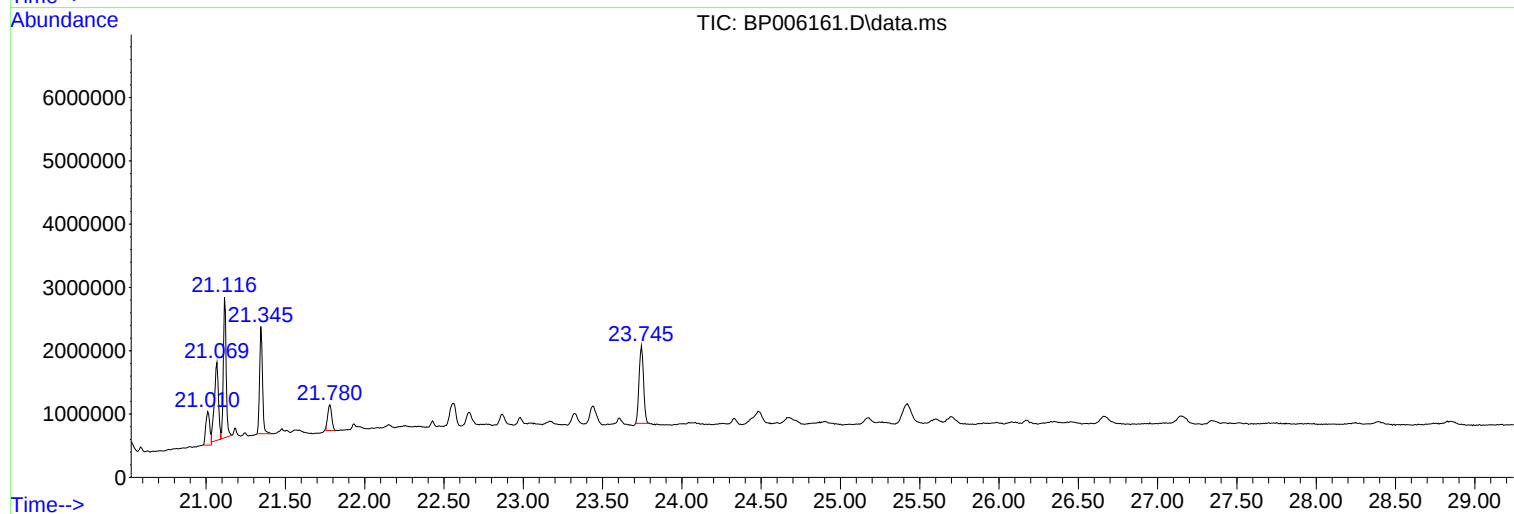
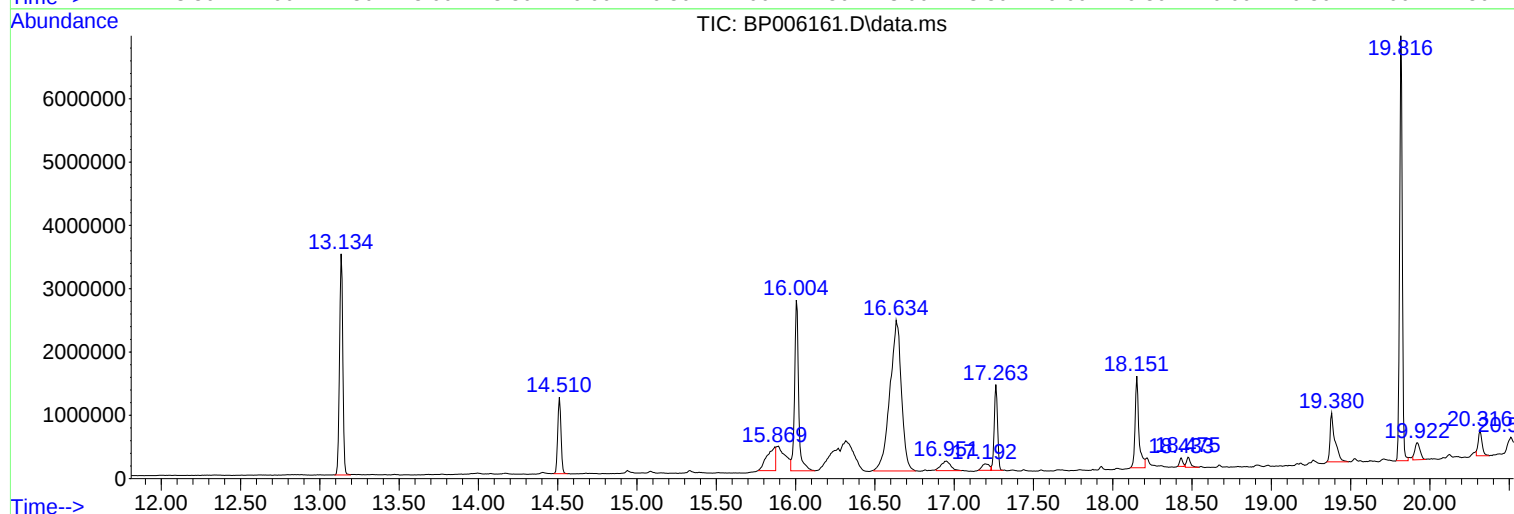
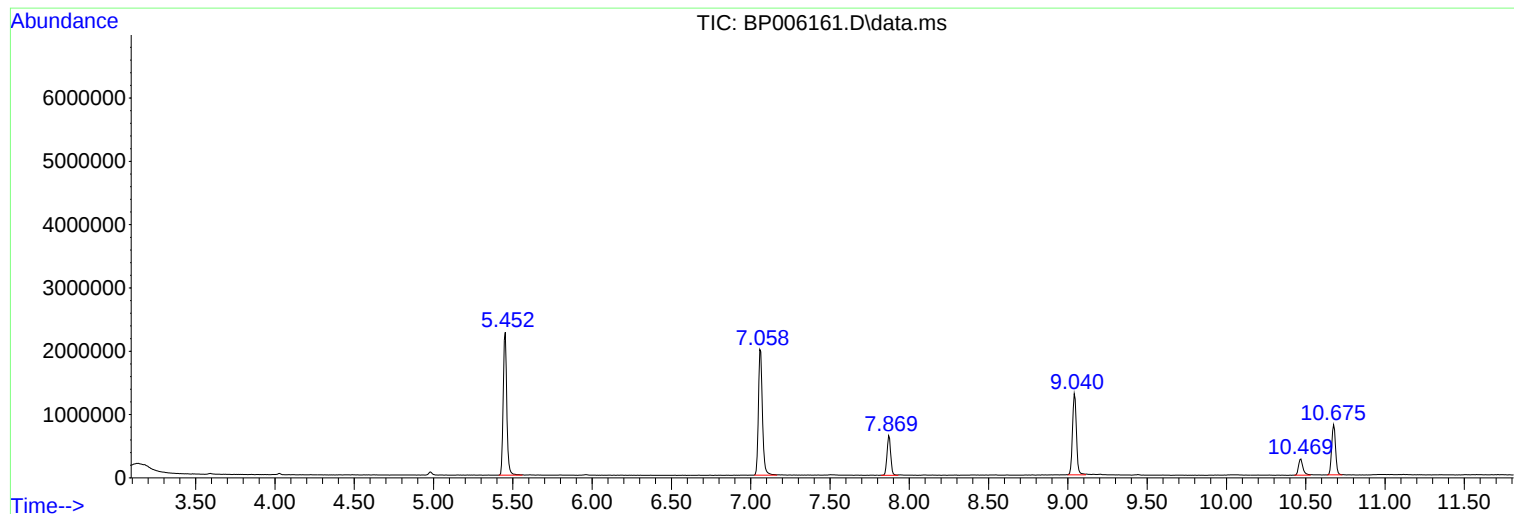
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP070721.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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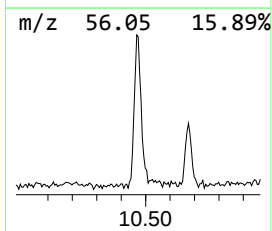
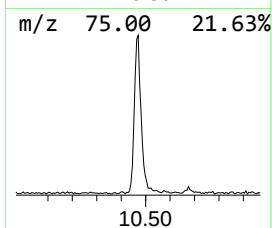
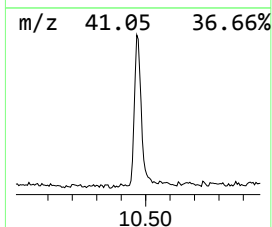
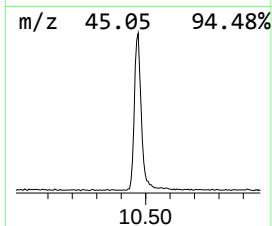
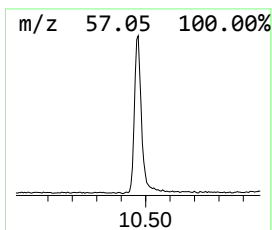
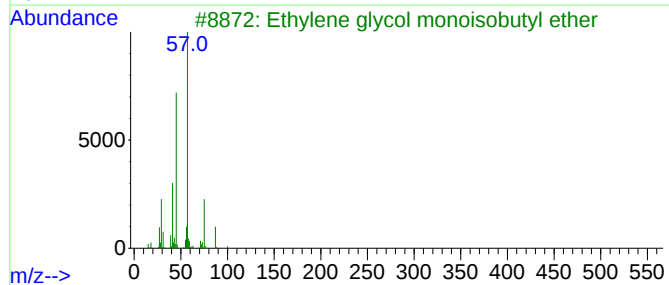
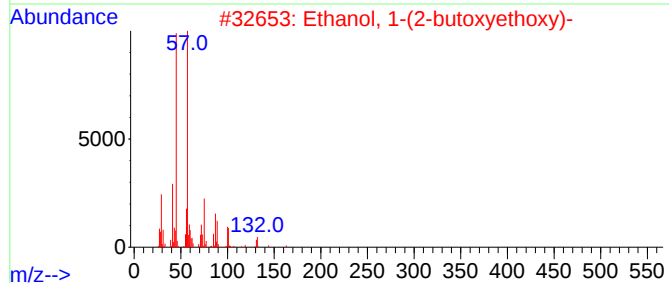
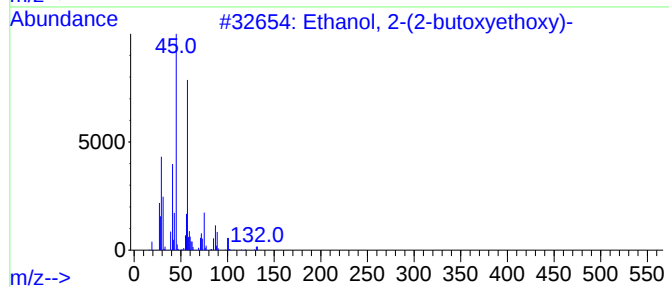
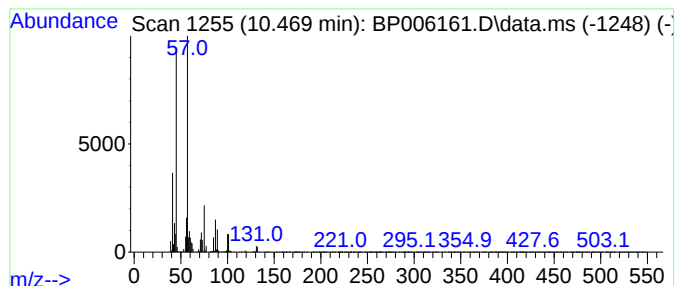
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Ethanol, 2-(2-butoxyethoxy)- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.469	7.35 ng	496296	Naphthalene-d8	10.675

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
2			Ethanol, 1-(2-butoxyethoxy)-	162	C8H18O3	054446-78-5	90
3			Ethylene glycol monoisobutyl ether	118	C6H14O2	004439-24-1	59
4			Ethanol, 2-[2-(2-butoxyethoxy)et...	206	C10H22O4	000143-22-6	59
5			3,3-Dimethylbutane-2-ol	102	C6H14O	000464-07-3	49



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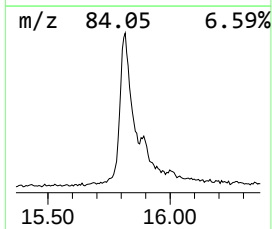
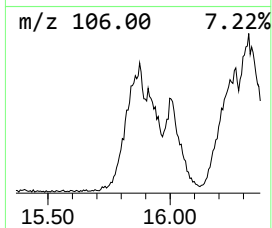
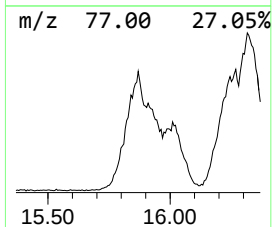
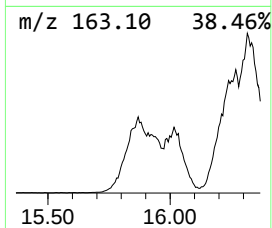
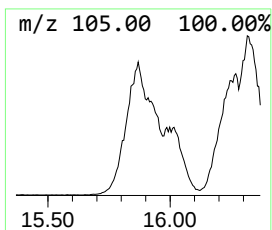
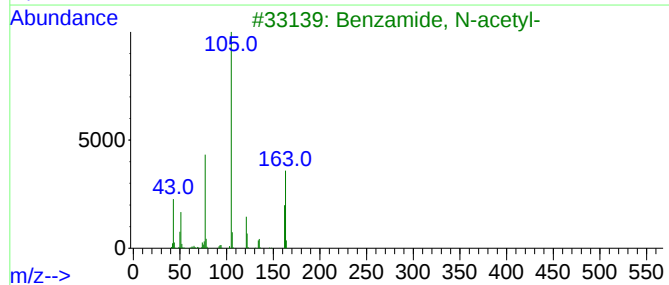
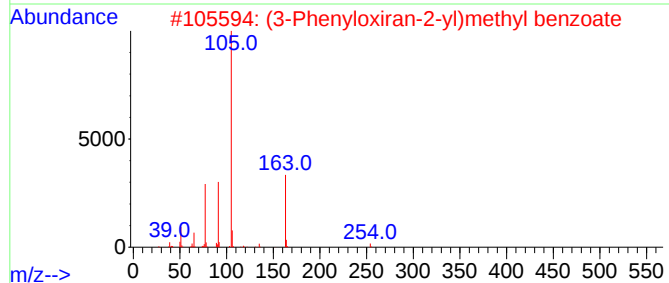
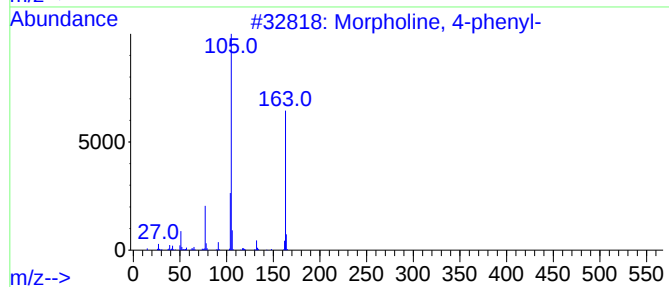
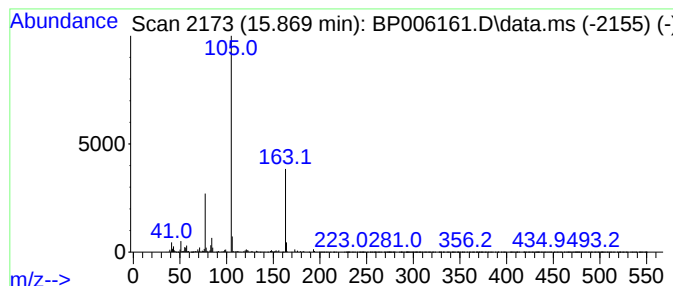
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Peak Number 2 unknown15.869 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.869	16.90 ng	1437060	Acenaphthene-d10	14.510

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Morpholine, 4-phenyl-	163	C10H13NO	000092-53-5	45
2			(3-Phenyloxiran-2-yl)methyl benz...	254	C16H14O3	1000366-96-1	40
3			Benzamide, N-acetyl-	163	C9H9NO2	001575-95-7	9
4			Benzoic acid, 2-benzyloxy-3-brom...	348	C17H17BrO3	1000191-33-3	9
5			Malonic acid, 2-chloropropyl pro...	222	C9H15ClO4	1000349-02-7	9



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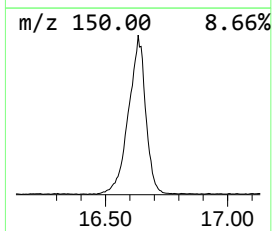
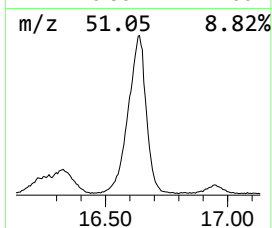
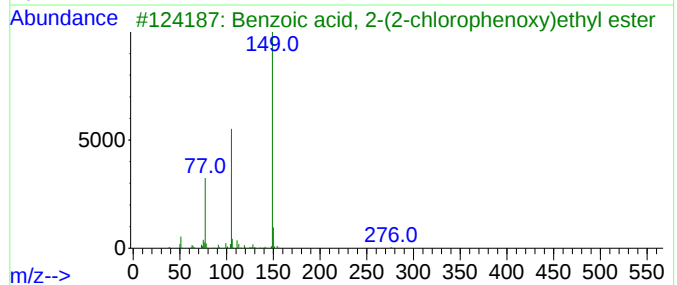
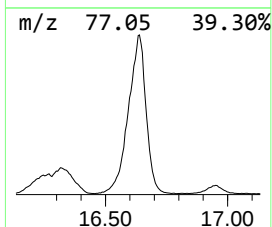
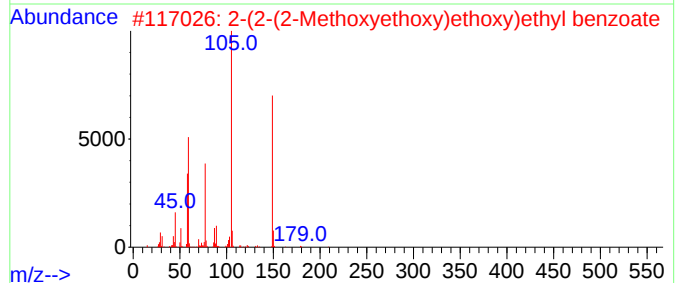
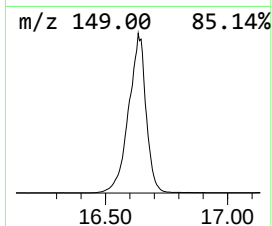
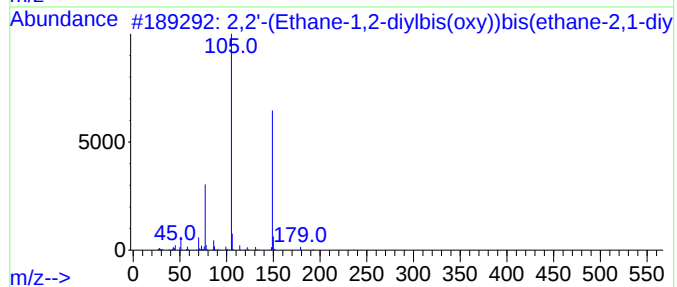
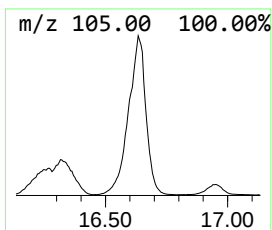
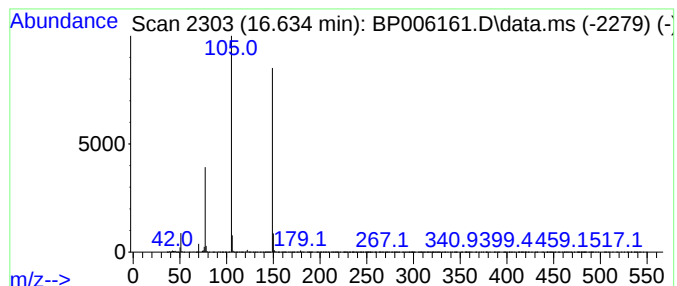
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 2,2'-(Ethane-1,2-diylbis(ox... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.	
16.634	121.25 ng	11998000	Phenanthrene-d10	17.263	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,2'-(Ethane-1,2-diylbis(oxy))bi...	358	C20H22O6	1000366-99-4	83
2	2-(2-(2-Methoxyethoxy)ethoxy)eth...	268	C14H20O5	1000366-99-1	74
3	Benzoic acid, 2-(2-chlorophenoxy...	276	C15H13ClO3	1000368-25-9	72
4	1,3-Dioxolane, 2-(methoxymethyl)...	194	C11H14O3	1000156-71-2	59
5	2-(2-Carboxyvinyl)pyridine, trans	149	C8H7NO2	007340-22-9	49



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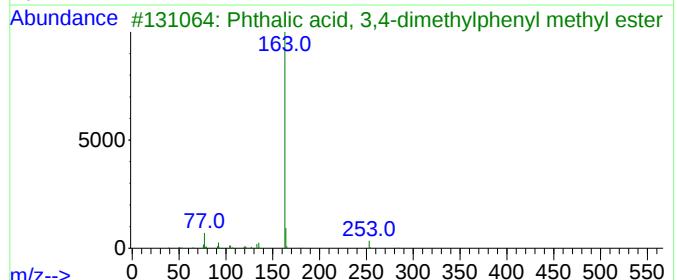
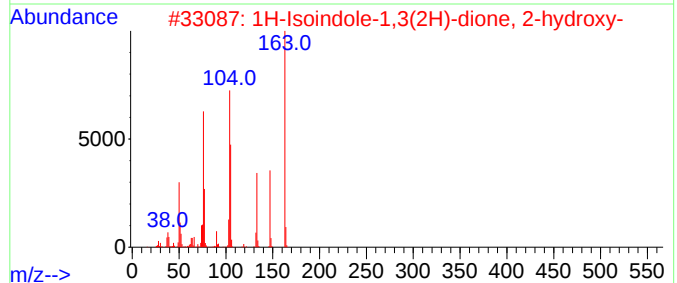
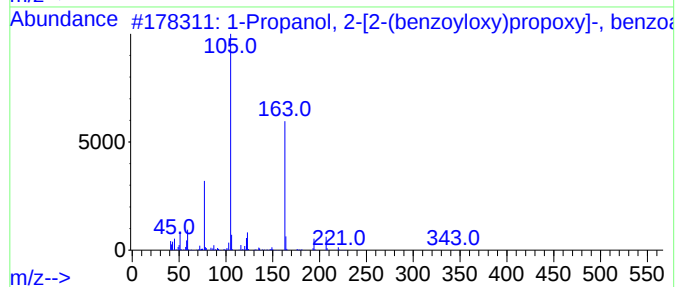
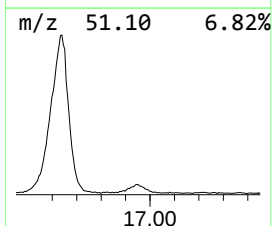
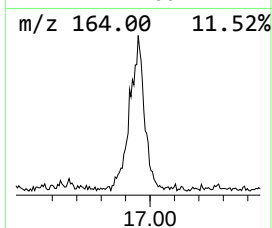
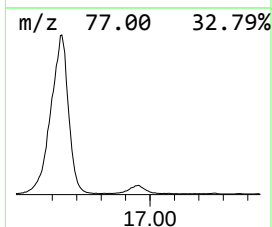
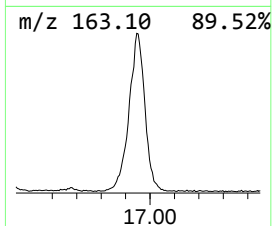
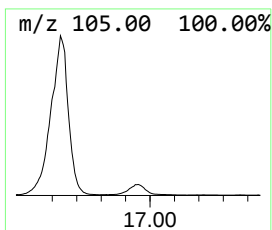
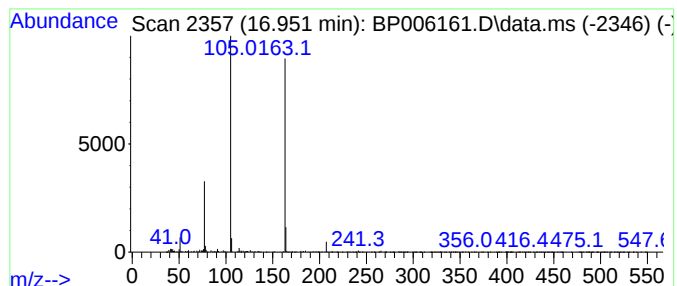
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Peak Number 4 1-Propanol, 2-[2-(benzoylox... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.951	6.39 ng	632673	Phenanthrene-d10	17.263

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Propanol, 2-[2-(benzoyloxy)pro...	342	C20H22O5	020109-39-1	64
2			1H-Isoindole-1,3(2H)-dione, 2-hy...	163	C8H5NO3	000524-38-9	50
3			Phthalic acid, 3,4-dimethylpheny...	284	C17H16O4	1000315-69-7	50
4			Glyoxylamide, N-phenyl-	163	C9H9NO2	032331-52-5	42
5			Morpholine, 4-phenyl-	163	C10H13NO	000092-53-5	40



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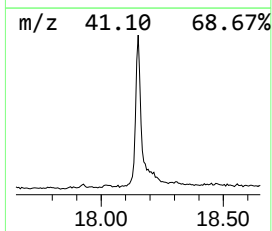
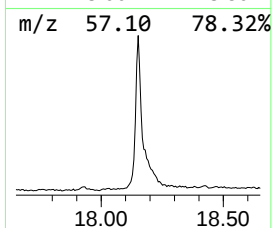
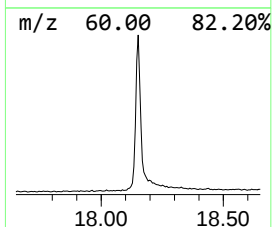
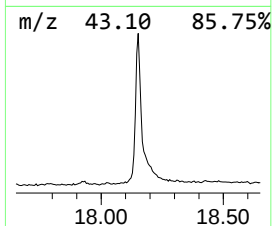
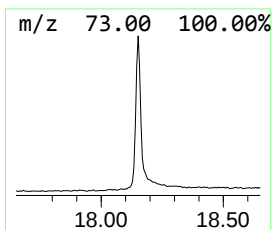
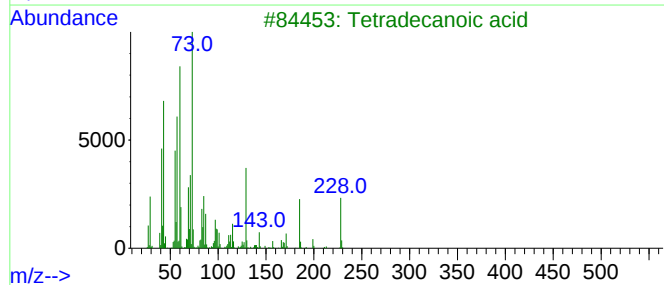
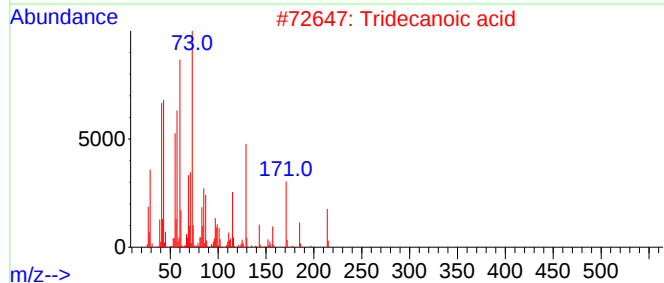
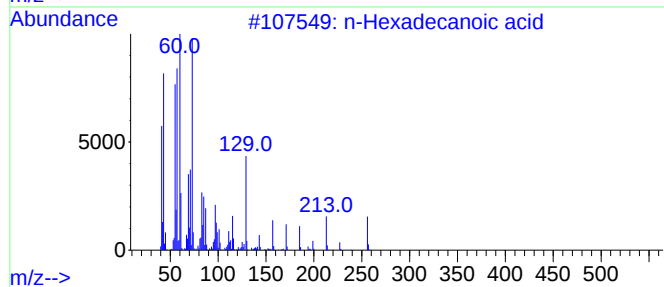
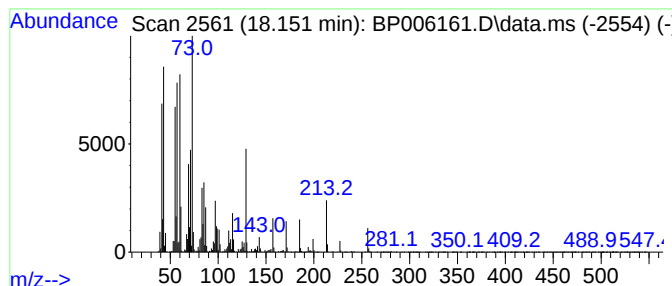
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 6 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.151	24.24 ng	2398160	Phenanthrene-d10	17.263

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	89
3			Tetradecanoic acid	228	C14H28O2	000544-63-8	83
4			Pentadecanoic acid	242	C15H30O2	001002-84-2	81
5			n-Decanoic acid	172	C10H20O2	000334-48-5	76



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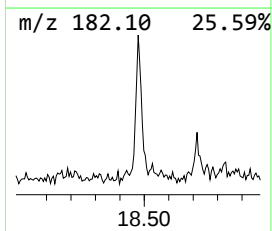
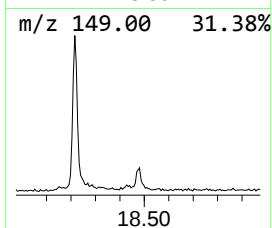
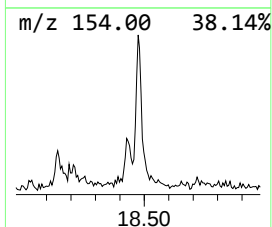
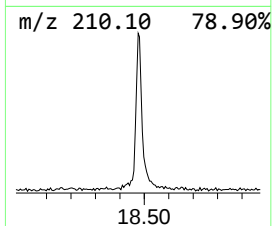
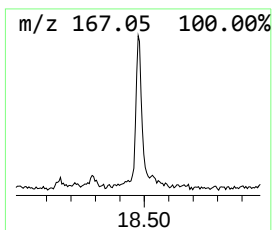
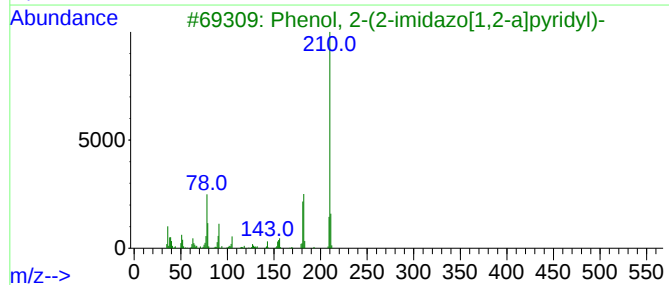
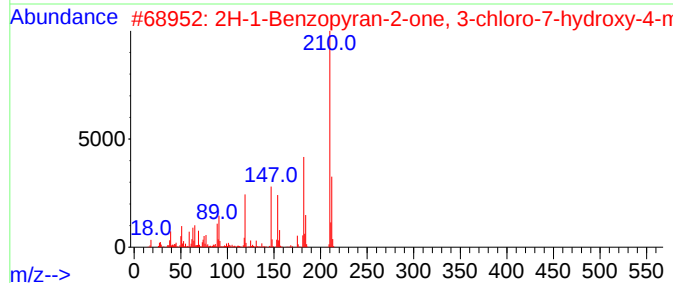
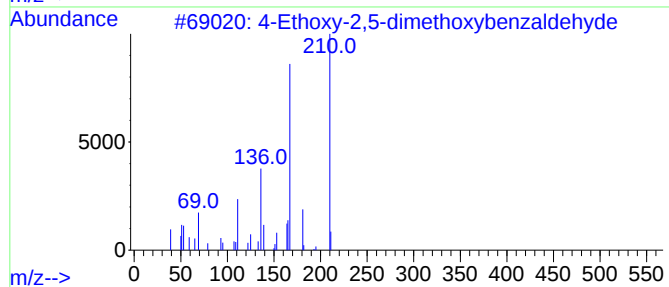
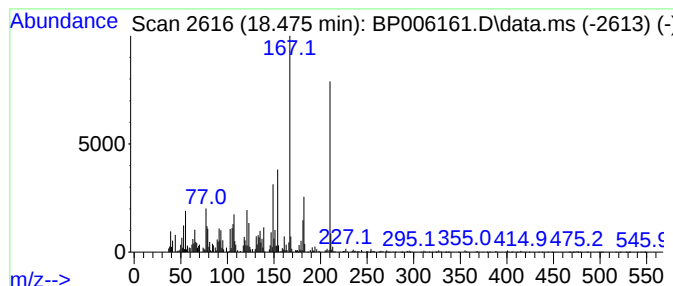
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 4-Ethoxy-2,5-dimethoxybenza... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.475	2.66 ng	262879	Phenanthrene-d10	17.263

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4-Ethoxy-2,5-dimethoxybenzaldehyde	210	C11H14O4	1000121-30-2	53
2			2H-1-Benzopyran-2-one, 3-chloro-...	210	C10H7ClO3	006174-86-3	38
3			Phenol, 2-(2-imidazo[1,2-a]pyrid...	210	C13H10N2O	057636-32-5	30
4			4-Chloro-1-methyl-1,5-naphthyrid...	210	C9H7ClN2O2	1000213-59-5	27
5			2H-1,3-Benzoxathiole, 5-formyl-7...	210	C9H6O4S	004991-69-9	27



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070921\
Data File : BP006161.D
Acq On : 09 Jul 2021 15:15
Operator : CG/JU
Sample : M2990-13
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
TP-J

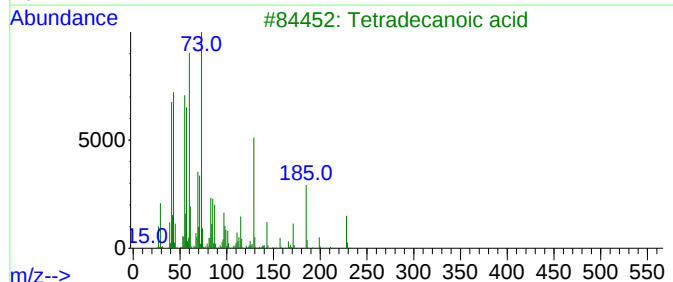
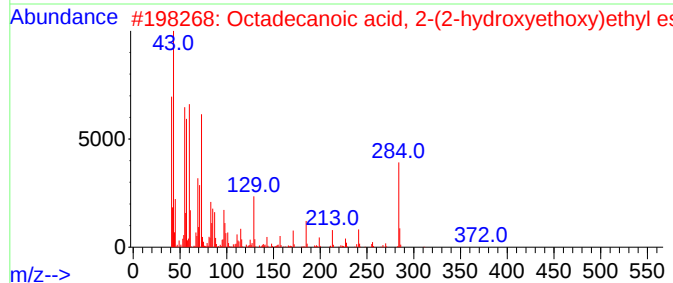
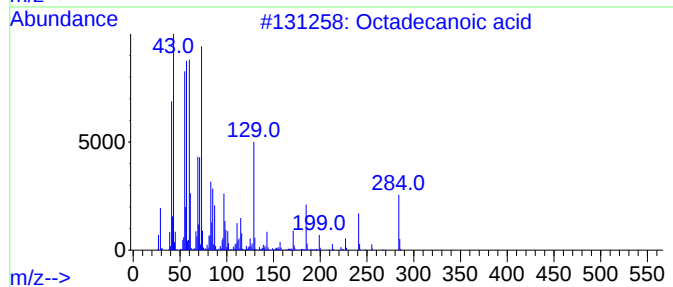
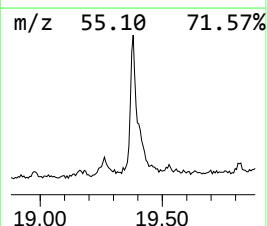
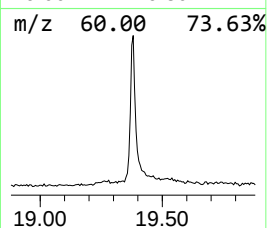
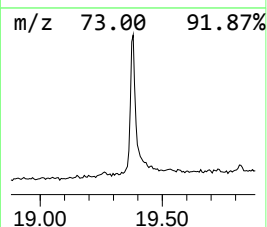
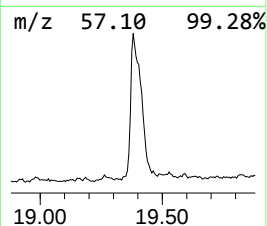
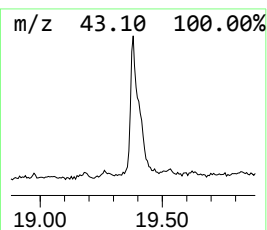
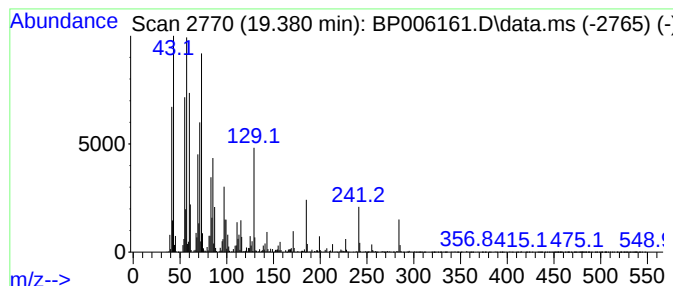
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 Octadecanoic acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.381	15.03 ng	1635470	Chrysene-d12	21.345

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid	284	C18H36O2	000057-11-4	99
2			Octadecanoic acid, 2-(2-hydroxye...	372	C22H44O4	000106-11-6	72
3			Tetradecanoic acid	228	C14H28O2	000544-63-8	70
4			Pentadecanoic acid	242	C15H30O2	001002-84-2	68
5			n-Decanoic acid	172	C10H20O2	000334-48-5	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070921\
Data File : BP006161.D
Acq On : 09 Jul 2021 15:15
Operator : CG/JU
Sample : M2990-13
Misc :
ALS Vial : 5 Sample Multiplier: 1

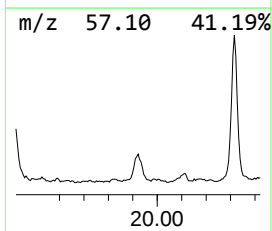
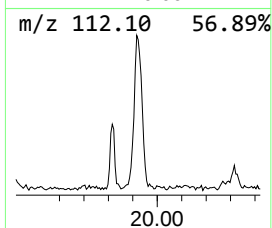
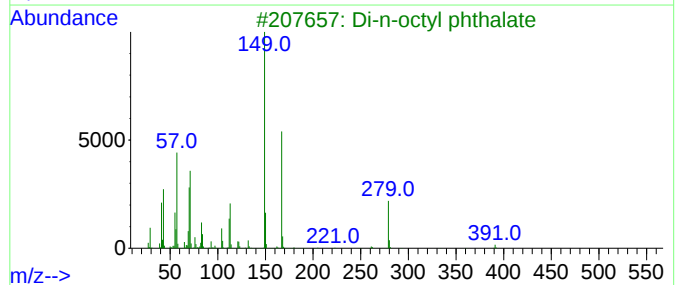
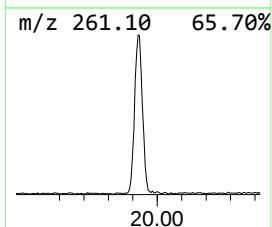
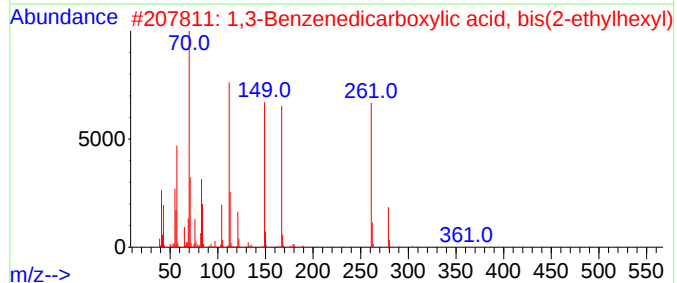
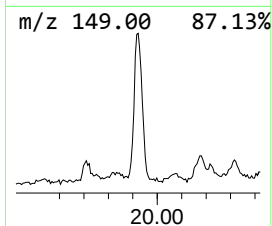
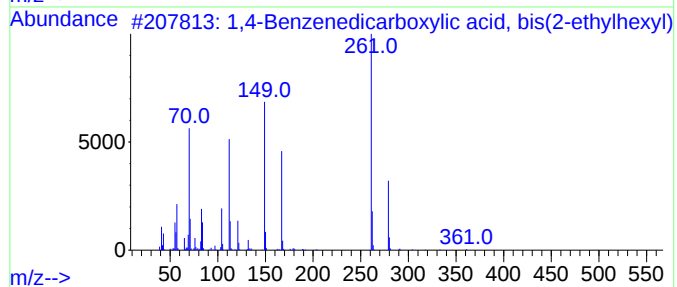
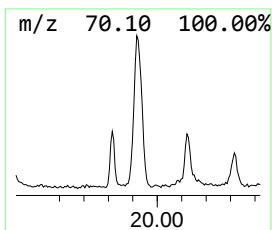
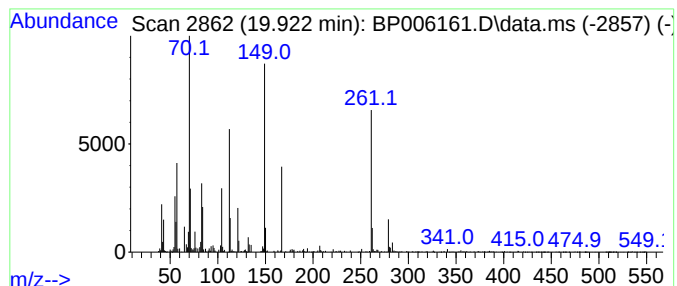
Instrument :
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ClientSampleId :
TP-J

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP070721.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 1,4-Benzenedicarboxylic aci... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD		R.T.	
19.922	5.62 ng	611606	Chrysene-d12		21.345	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	1,4-Benzenedicarboxylic acid, bi...		390	C24H38O4	006422-86-2	72
2	1,3-Benzenedicarboxylic acid, bi...		390	C24H38O4	000137-89-3	53
3	Di-n-octyl phthalate		390	C24H38O4	000117-84-0	46
4	Terephthalic acid, monochloride,...		296	C16H21ClO3	1000324-22-9	38
5	Diisooctyl phthalate		390	C24H38O4	000131-20-4	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070921\
Data File : BP006161.D
Acq On : 09 Jul 2021 15:15
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Sample : M2990-13
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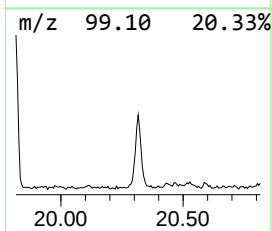
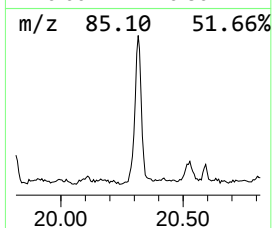
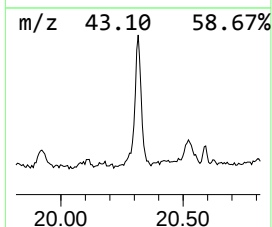
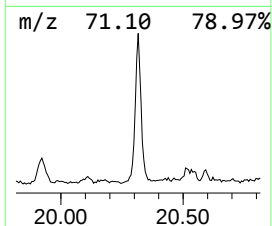
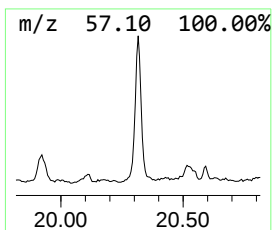
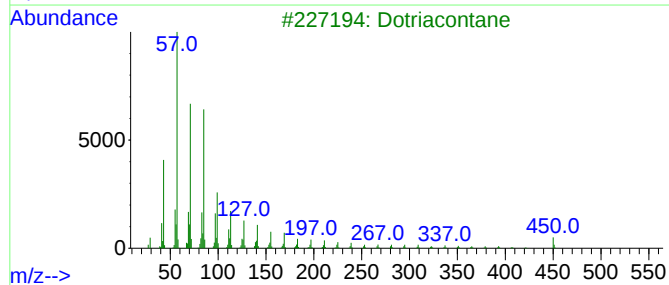
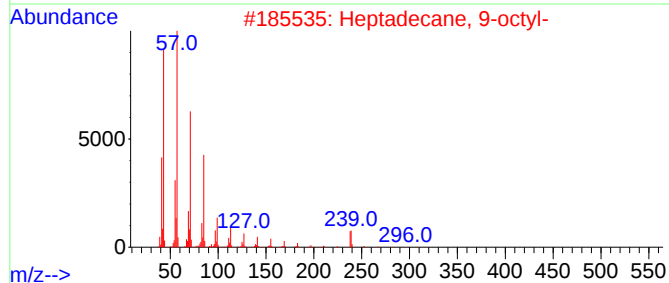
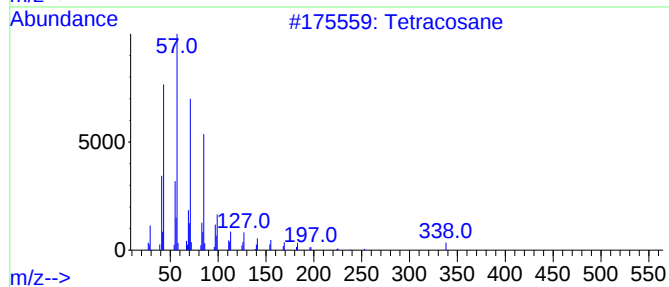
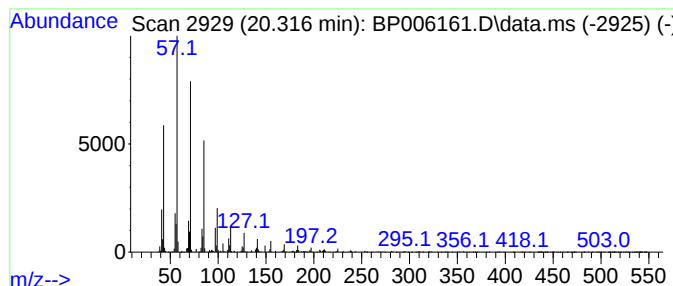
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP070721.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 Tetracosane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.	
20.316	6.21 ng	675632	Chrysene-d12	21.345	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetracosane	338	C24H50	000646-31-1	93
2	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	93
3	Dotriacontane	451	C32H66	000544-85-4	93
4	Eicosane, 9-octyl-	394	C28H58	013475-77-9	93
5	Heneicosane	296	C21H44	000629-94-7	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070921\
Data File : BP006161.D
Acq On : 09 Jul 2021 15:15
Operator : CG/JU
Sample : M2990-13
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
TP-J

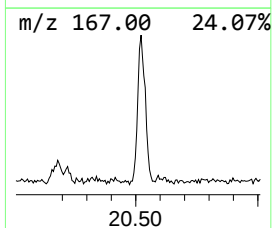
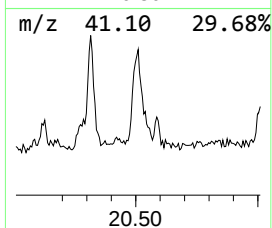
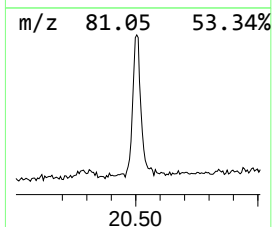
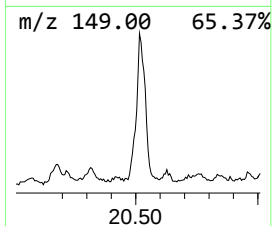
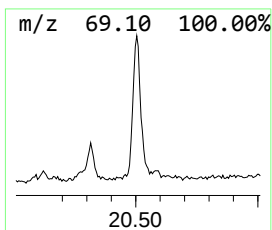
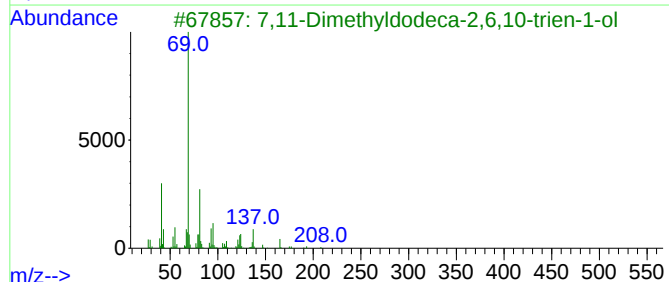
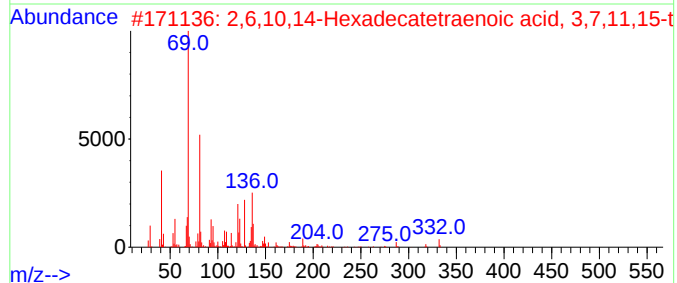
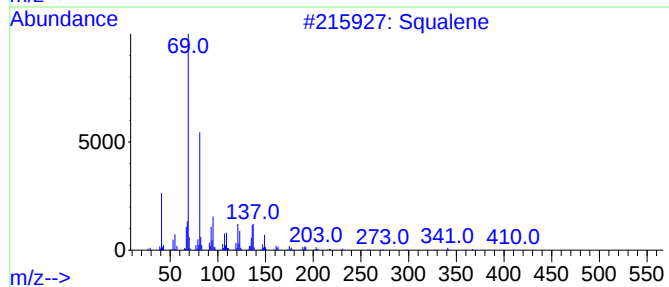
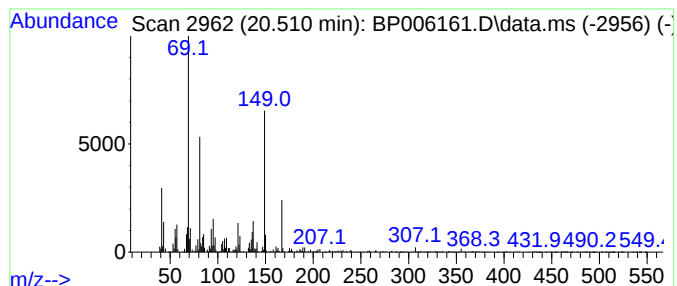
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 Squalene Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.510	6.10 ng	663473	Chrysene-d12	21.345

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Squalene	410	C30H50	000111-02-4	64
2			2,6,10,14-Hexadecatetraenoic aci...	332	C22H36O2	024035-35-6	47
3			7,11-Dimethyldodeca-2,6,10-trien...	208	C14H24O	125529-10-4	43
4			2,6-Octadiene, 4-methyl-	124	C9H16	074498-94-5	35
5			2-((4-Methylpentan-2-yloxy)carbo...	250	C14H18O4	1000373-82-5	14



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070921\
Data File : BP006161.D
Acq On : 09 Jul 2021 15:15
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Sample : M2990-13
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
TP-J

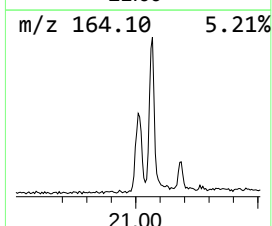
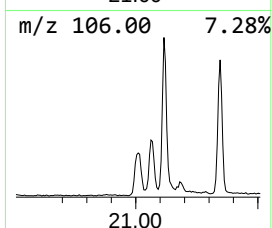
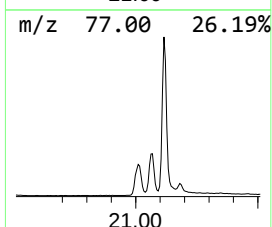
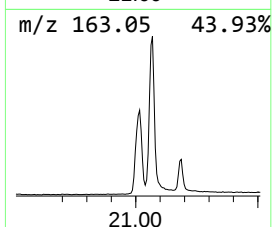
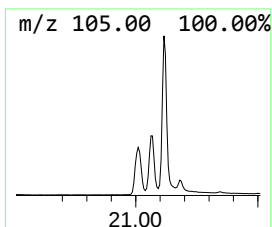
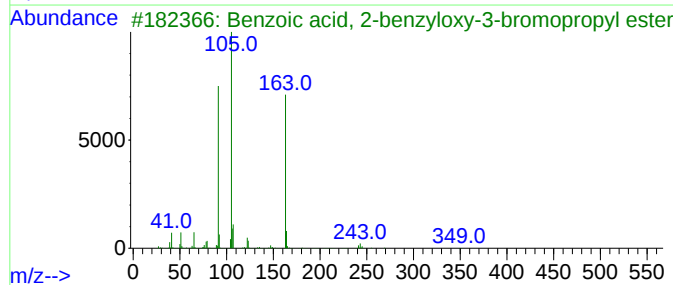
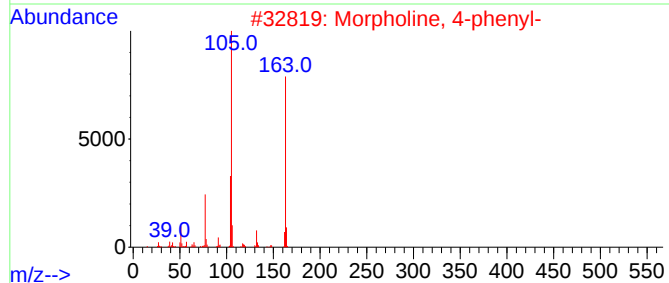
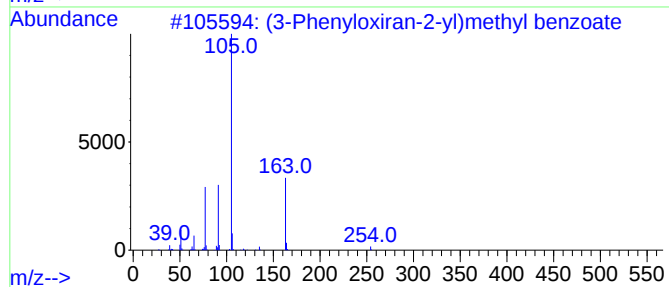
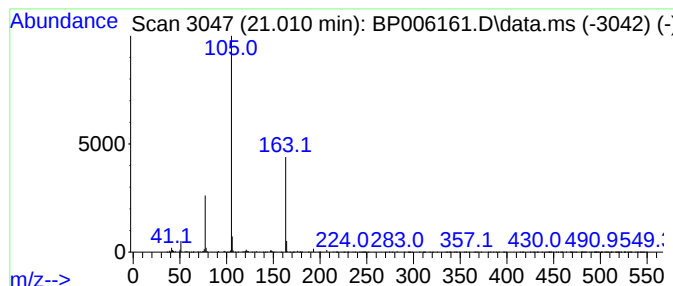
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 12 (3-Phenyloxiran-2-yl)methyl... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.010	8.54 ng	929941	Chrysene-d12	21.345

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			(3-Phenyloxiran-2-yl)methyl benz...	254	C16H14O3	1000366-96-1	56
2			Morpholine, 4-phenyl-	163	C10H13NO	000092-53-5	53
3			Benzoic acid, 2-benzyloxy-3-brom...	348	C17H17BrO3	1000191-33-3	33
4			Benzamide, N-acetyl-	163	C9H9NO2	001575-95-7	9
5			1-Benzoyl-2-trichloroacetylhydra...	280	C9H7Cl3N2O2	090273-66-8	9



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070921\
Data File : BP006161.D
Acq On : 09 Jul 2021 15:15
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Sample : M2990-13
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
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ClientSampleId :
TP-J

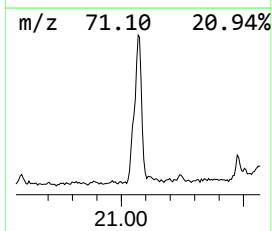
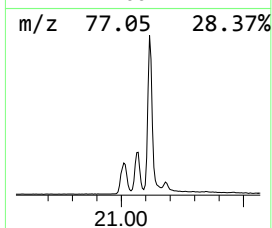
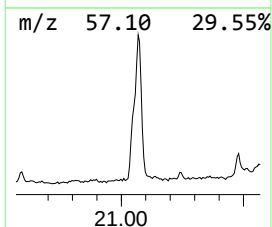
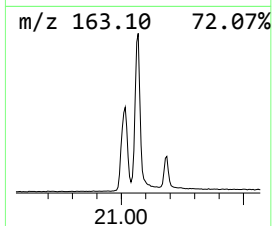
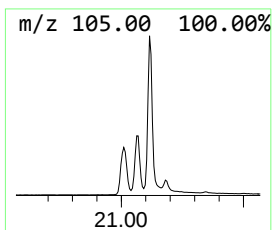
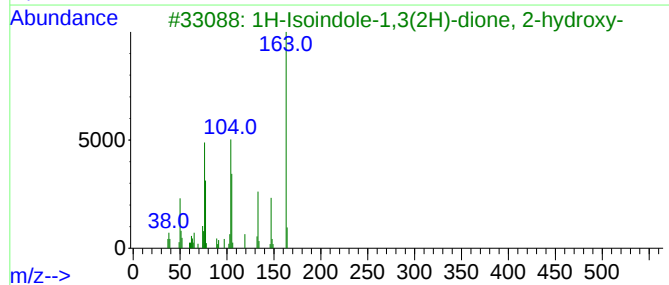
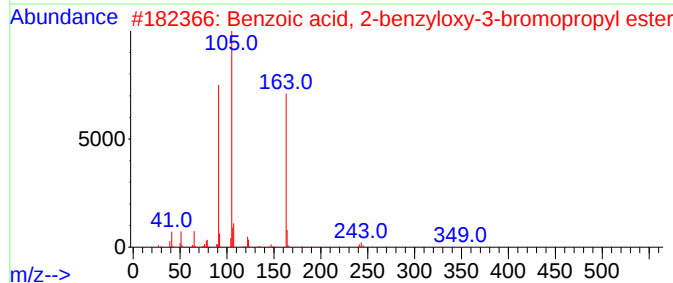
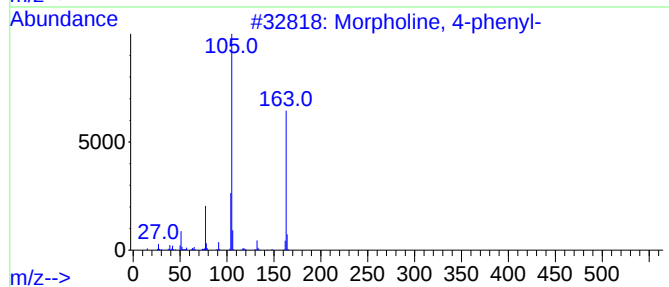
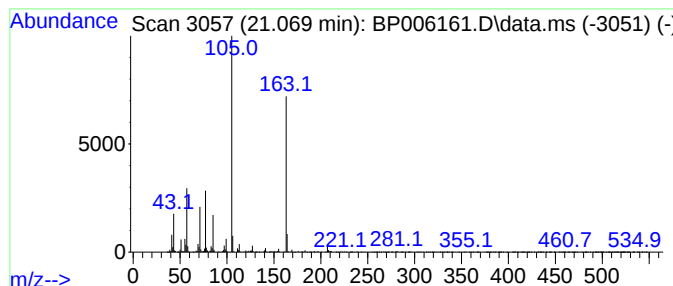
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 13 unknown21.069 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.069	20.28 ng	2207030	Chrysene-d12	21.345

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Morpholine, 4-phenyl-	163	C10H13NO	000092-53-5	40
2			Benzoic acid, 2-benzyloxy-3-brom...	348	C17H17BrO3	1000191-33-3	38
3			1H-Isoindole-1,3(2H)-dione, 2-hy...	163	C8H5NO3	000524-38-9	38
4			Ethanone, 1-[2-(dimethylamino)ph...	163	C10H13NO	010336-55-7	38
5			Glyoxylamide, N-phenyl-	163	C9H9NO2	032331-52-5	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070921\
Data File : BP006161.D
Acq On : 09 Jul 2021 15:15
Operator : CG/JU
Sample : M2990-13
Misc :
ALS Vial : 5 Sample Multiplier: 1

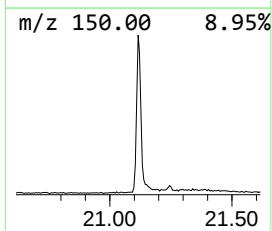
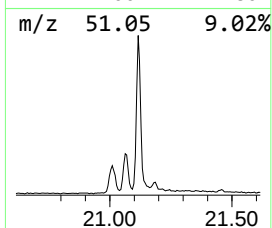
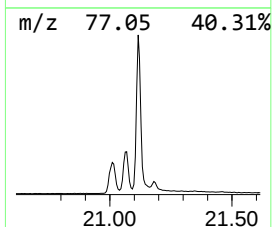
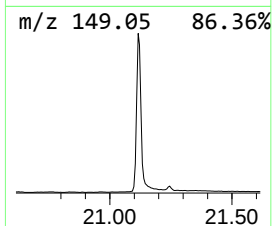
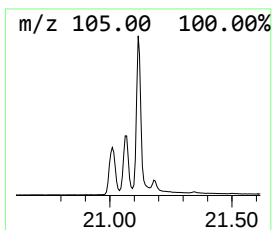
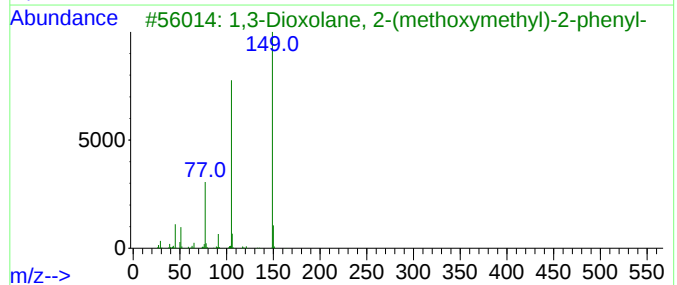
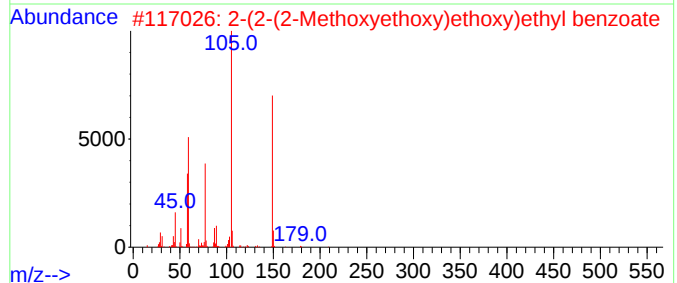
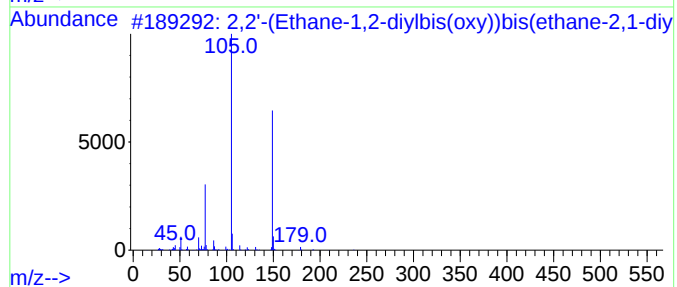
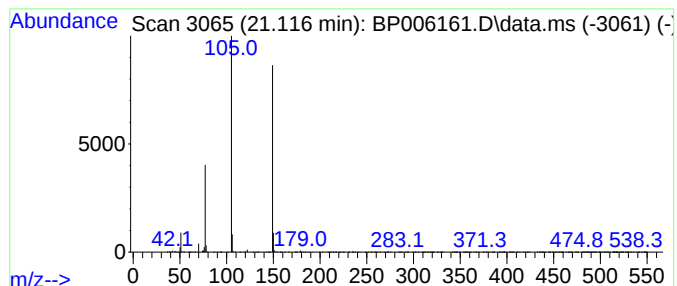
Instrument :
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ClientSampleId :
TP-J

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP070721.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 14 2-(2-(2-Methoxyethoxy)ethox... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.		
21.116	25.15 ng	2737130	Chrysene-d12	21.345		
Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,2'	-(Ethane-1,2-diylbis(oxy))bi...	358	C20H22O6	1000366-99-4	83
2	2-(2-(2-Methoxyethoxy)ethoxy)eth...	268	C14H20O5	1000366-99-1	74	
3	1,3-Dioxolane, 2-(methoxymethyl)...	194	C11H14O3	1000156-71-2	59	
4	3,6,9,12-Tetraoxatetradecane-1,1...	446	C24H30O8	1000366-97-0	50	
5	2-(2-Carboxyvinyl)pyridine, trans	149	C8H7NO2	007340-22-9	49	



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070921\
Data File : BP006161.D
Acq On : 09 Jul 2021 15:15
Operator : CG/JU
Sample : M2990-13
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
TP-J

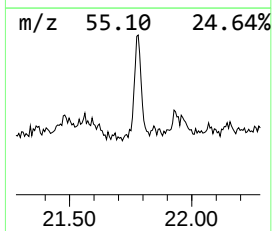
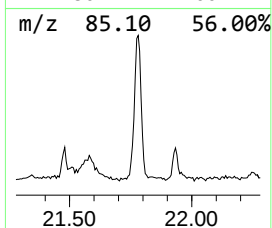
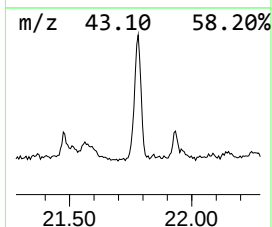
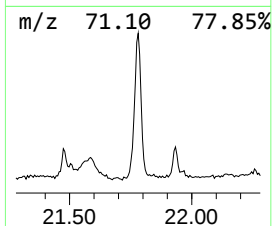
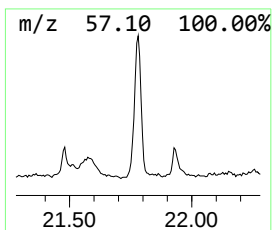
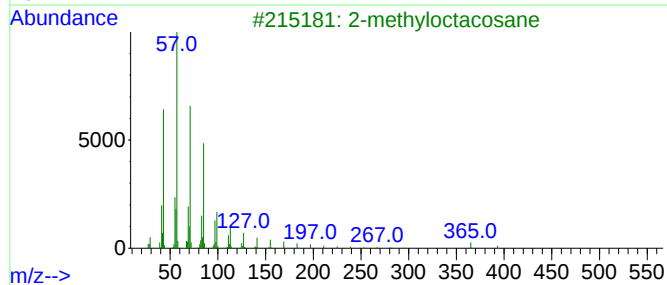
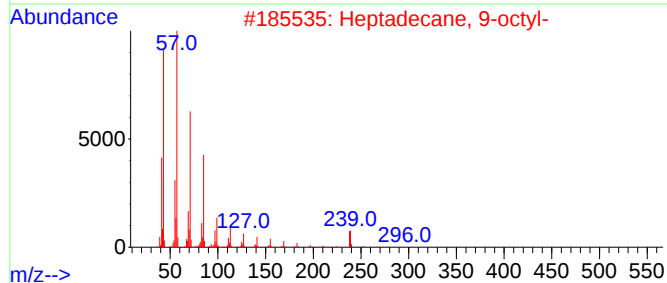
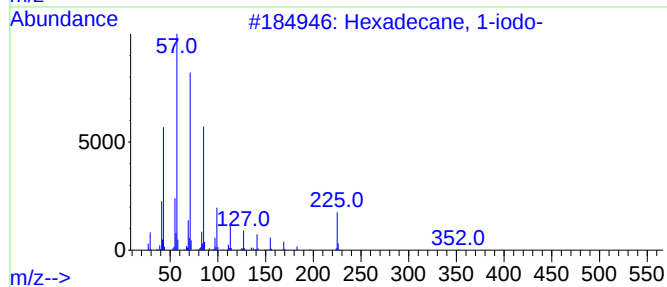
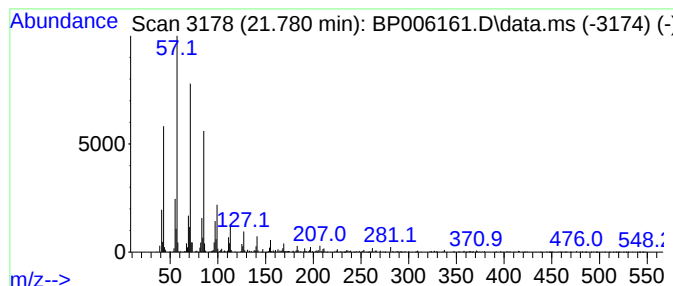
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 15 Hexadecane, 1-iodo- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.780	6.36 ng	692103	Chrysene-d12	21.345

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexadecane, 1-iodo-	352	C16H33I	000544-77-4	95
2			Heptadecane, 9-octyl-	352	C25H52	007225-64-1	93
3			2-methyloctacosane	408	C29H60	1000376-72-8	91
4			Heneicosane	296	C21H44	000629-94-7	91
5			Octacosane	394	C28H58	000630-02-4	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070921\
Data File : BP006161.D
Acq On : 09 Jul 2021 15:15
Operator : CG/JU
Sample : M2990-13
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
TP-J

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP070721.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Ethanol, 2-(2-b...	10.469	7.3	ng	496296	2	10.675	1350580	20.0
unknown15.869	15.869	16.9	ng	1437060	3	14.510	1700270	20.0
2,2'-(Ethane-1,...	16.634	121.3	ng	11998000	4	17.263	1979010	20.0
1-Propanol, 2-[...	16.951	6.4	ng	632673	4	17.263	1979010	20.0
n-Hexadecanoic ...	18.151	24.2	ng	2398160	4	17.263	1979010	20.0
4-Ethoxy-2,5-di...	18.475	2.7	ng	262879	4	17.263	1979010	20.0
Octadecanoic acid	19.381	15.0	ng	1635470	5	21.345	2176740	20.0
1,4-Benzenedica...	19.922	5.6	ng	611606	5	21.345	2176740	20.0
Tetracosane	20.316	6.2	ng	675632	5	21.345	2176740	20.0
Squalene	20.510	6.1	ng	663473	5	21.345	2176740	20.0
(3-Phenyloxiran...	21.010	8.5	ng	929941	5	21.345	2176740	20.0
unknown21.069	21.069	20.3	ng	2207030	5	21.345	2176740	20.0
2-(2-(2-Methoxy...	21.116	25.1	ng	2737130	5	21.345	2176740	20.0
Hexadecane, 1-i...	21.780	6.4	ng	692103	5	21.345	2176740	20.0