

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071221\
 Data File : BP006188.D
 Acq On : 12 Jul 2021 19:04
 Operator : CG/JU
 Sample : M2997-01
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 RBR-200029

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: BP006188.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.446	395	401	427	rBV	1376770	2408894	22.28%	4.397%
2	7.057	669	675	691	rBV	1145329	2280856	21.09%	4.163%
3	7.722	777	788	795	rBV2	165191	466695	4.32%	0.852%
4	7.863	803	812	818	rBV	673111	1140224	10.55%	2.081%
5	7.940	818	825	867	rBV	4292759	10812522	100.00%	19.737%
6	9.028	1004	1010	1019	rBV	859312	1544582	14.29%	2.819%
7	9.228	1034	1044	1050	rBV4	191021	610860	5.65%	1.115%
8	9.351	1062	1065	1081	rVB8	219993	621223	5.75%	1.134%
9	10.251	1207	1218	1222	rBV5	382460	1045735	9.67%	1.909%
10	10.451	1246	1252	1257	rBV	604220	1074518	9.94%	1.961%
11	10.510	1257	1262	1274	rVB6	565293	1139744	10.54%	2.080%
12	10.663	1283	1288	1297	rVB	899301	1623203	15.01%	2.963%
13	13.122	1700	1706	1714	rVB	2669985	3696649	34.19%	6.748%
14	14.151	1876	1881	1889	rBV	5738686	7808580	72.22%	14.253%
15	14.222	1890	1893	1899	rVV3	300801	544653	5.04%	0.994%
16	14.322	1906	1910	1916	rVB	534920	803966	7.44%	1.468%
17	14.498	1936	1940	1947	rVB	1374909	1977086	18.29%	3.609%
18	15.275	2069	2072	2076	rBV	466331	624213	5.77%	1.139%
19	16.004	2192	2196	2207	rBV	772305	1939685	17.94%	3.541%
20	16.792	2325	2330	2346	rBV	2334696	7387568	68.32%	13.485%
21	17.063	2371	2376	2390	rBV	1819591	5232568	48.39%	9.551%

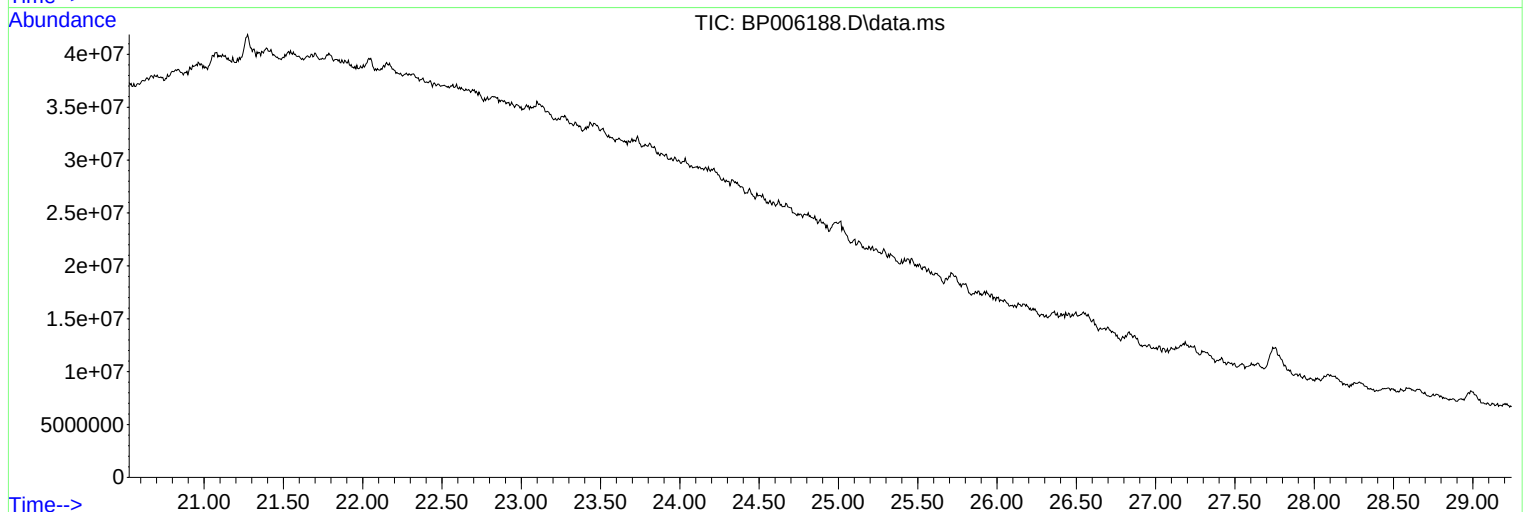
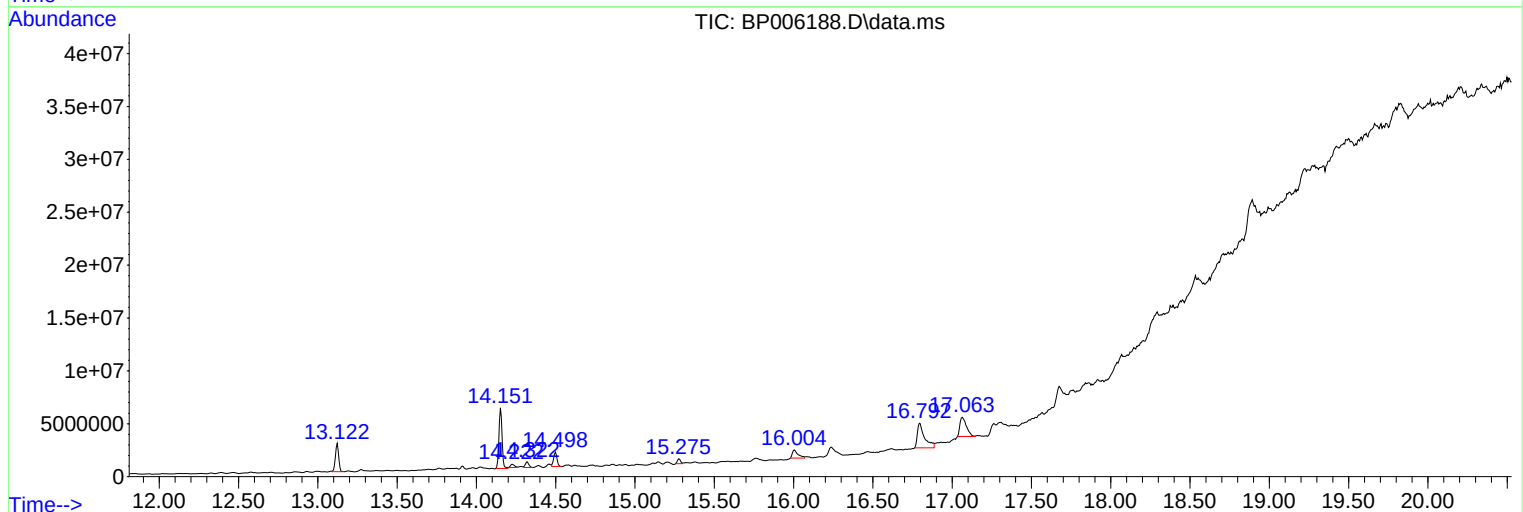
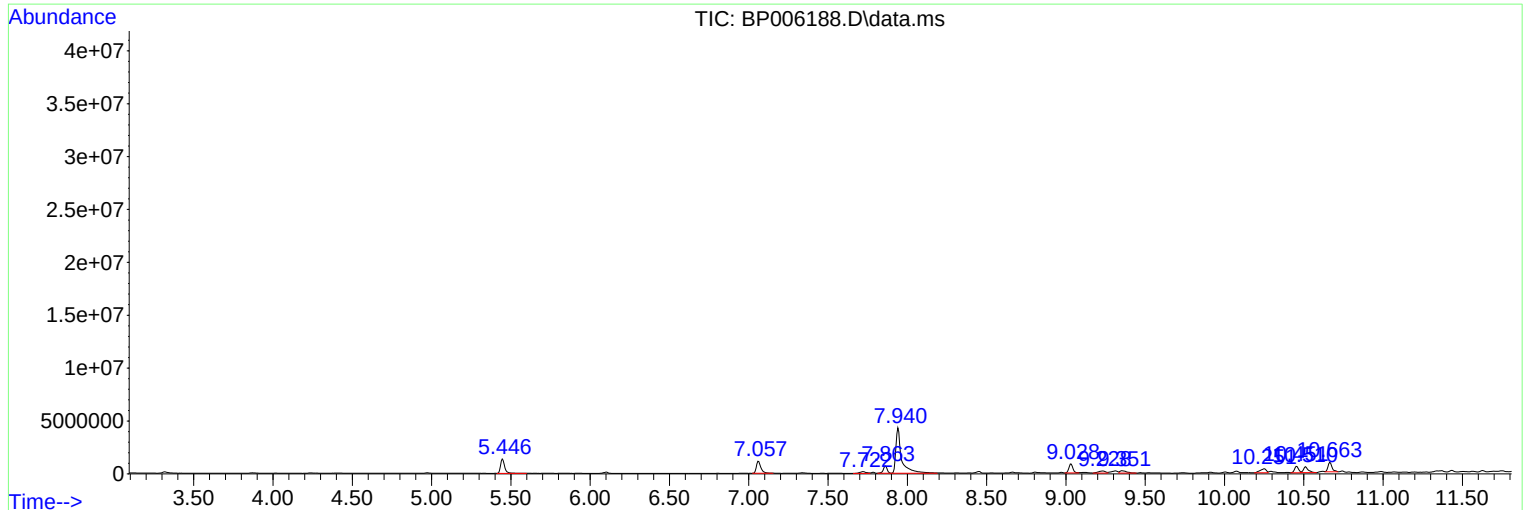
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP070721.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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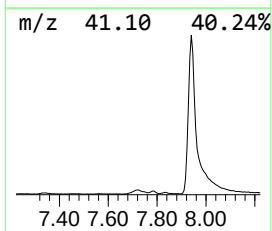
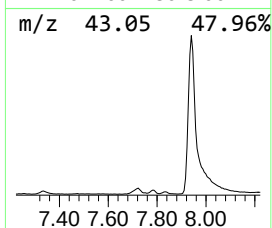
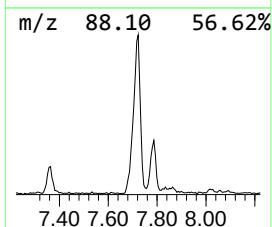
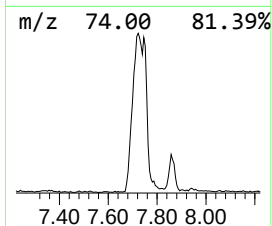
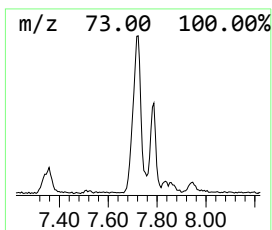
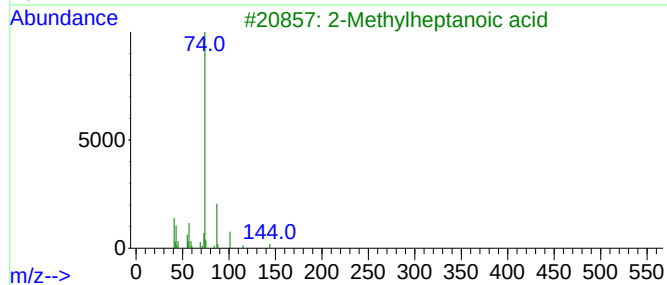
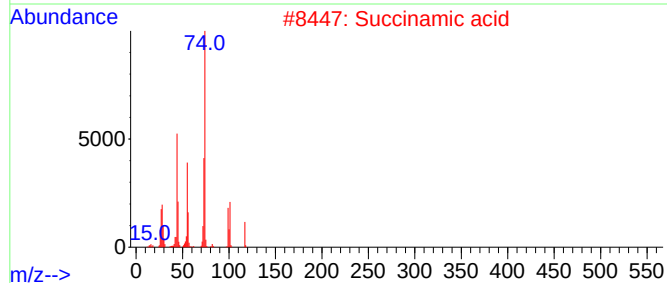
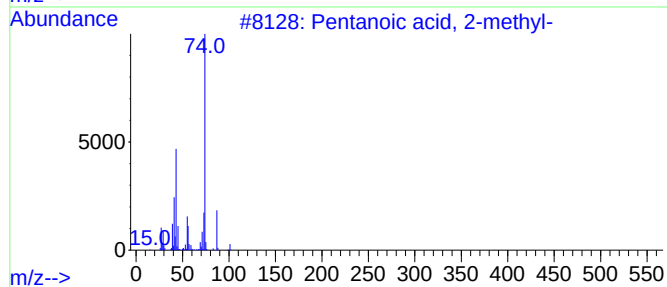
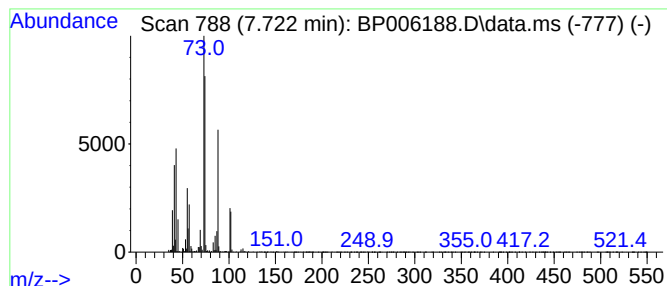
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 unknown7.722 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.722	8.19 ng	466695	1,4-Dichlorobenzene-d4	7.863

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentanoic acid, 2-methyl-	116	C6H12O2	000097-61-0	43
2			Succinamic acid	117	C4H7NO3	000638-32-4	38
3			2-Methylheptanoic acid	144	C8H16O2	001188-02-9	38
4			Pentanoic acid, 2-methyl-, butyl...	172	C10H20O2	006297-41-2	32
5			Butanoic acid, 2-methyl-	102	C5H10O2	000116-53-0	32



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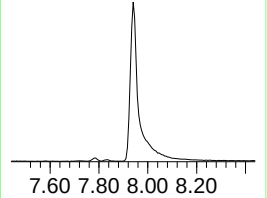
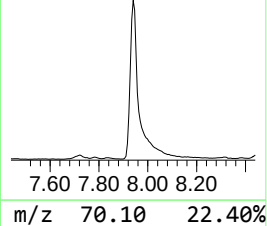
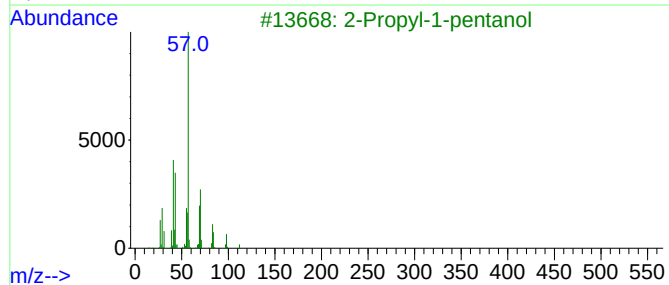
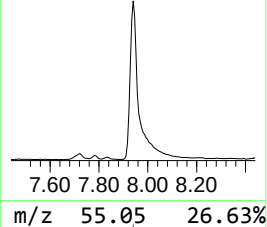
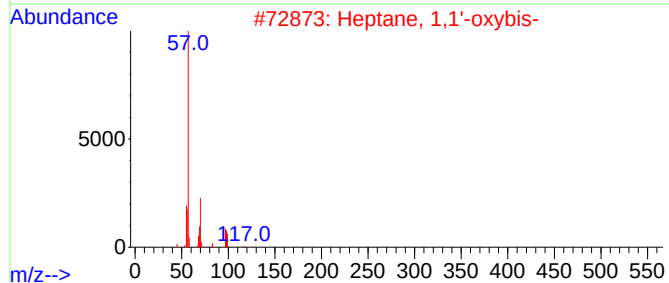
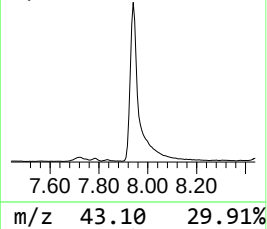
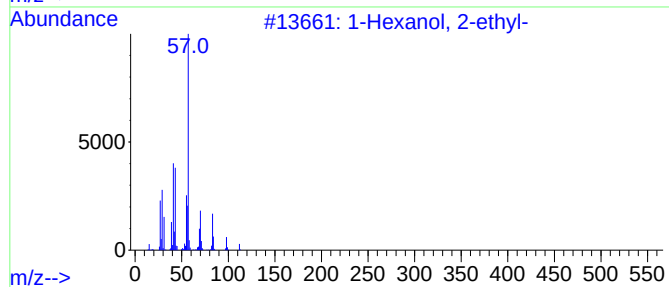
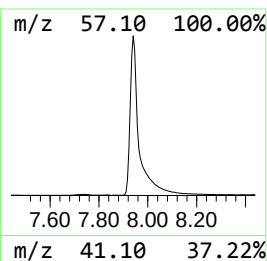
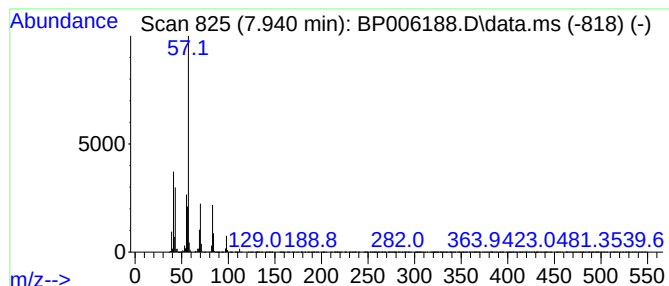
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 2 1-Hexanol, 2-ethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.940	189.66 ng	10812500	1,4-Dichlorobenzene-d4	7.863

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Hexanol, 2-ethyl-			130	C8H18O	000104-76-7	74
2	Heptane, 1,1'-oxybis-			214	C14H30O	000629-64-1	59
3	2-Propyl-1-pentanol			130	C8H18O	058175-57-8	53
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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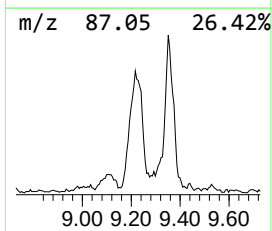
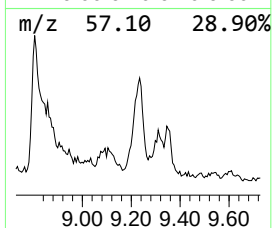
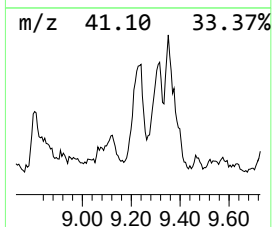
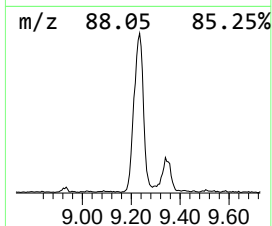
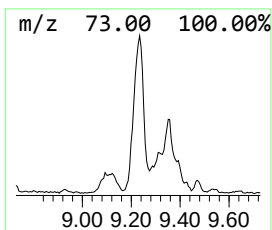
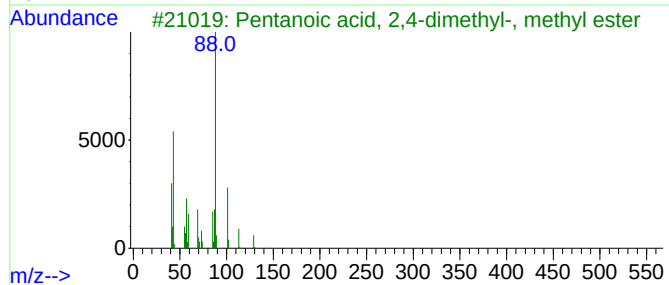
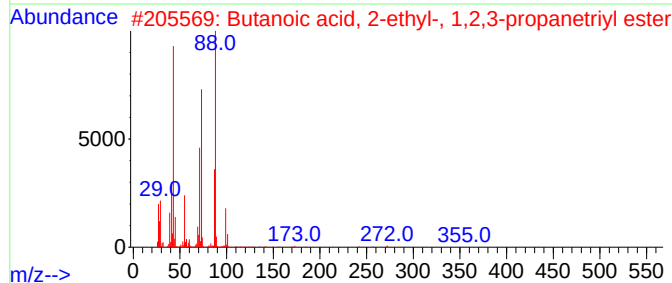
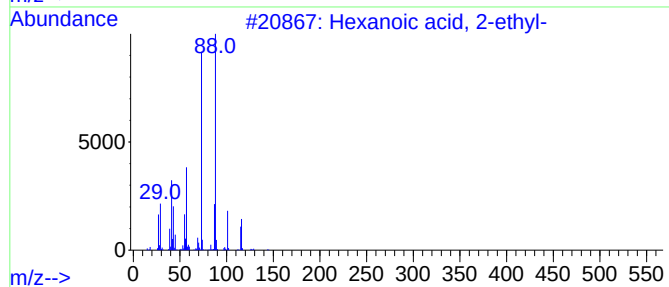
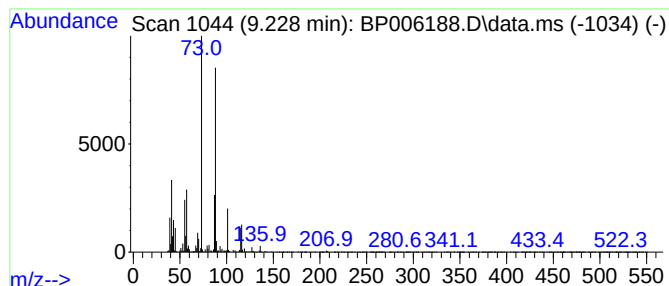
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Hexanoic acid, 2-ethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.228	10.71 ng	610860	1,4-Dichlorobenzene-d4	7.863

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	72
2			Butanoic acid, 2-ethyl-, 1,2,3-p...	386	C21H38O6	056554-54-2	53
3			Pentanoic acid, 2,4-dimethyl-, m...	144	C8H16O2	071672-33-8	50
4			1,4-Dioxane	88	C4H8O2	000123-91-1	43
5			Pentanoic acid, 2-methyl-, methy...	130	C7H14O2	002177-77-7	40



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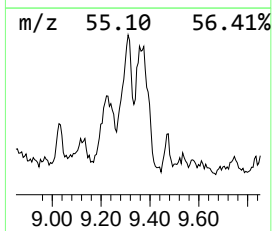
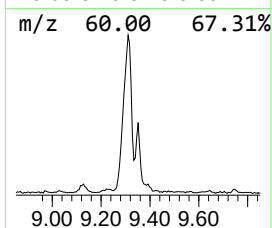
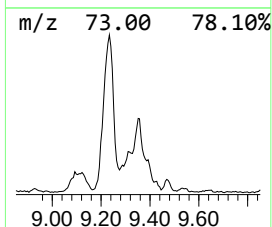
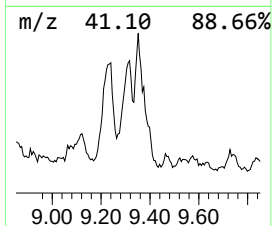
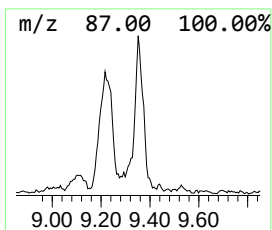
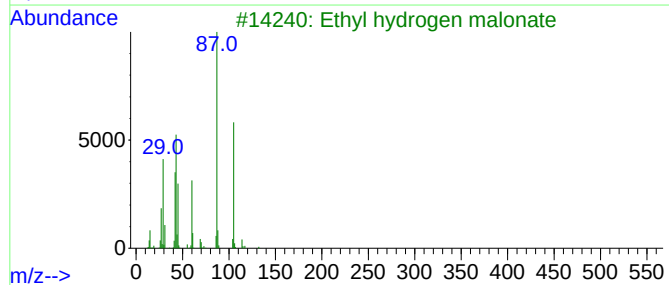
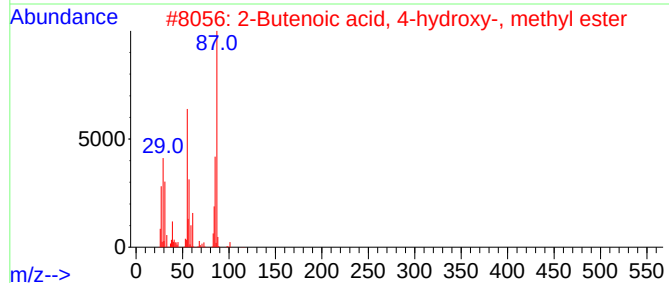
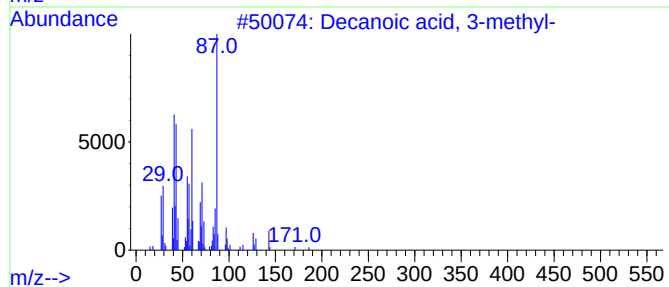
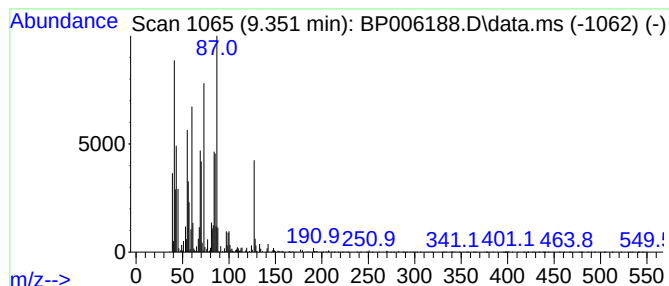
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TIC Integration Parameters: LSCINT.P

Peak Number 4 unknown9.351 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.351	7.65 ng	621223	Naphthalene-d8	10.663

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Decanoic acid, 3-methyl-	186	C11H22O2	060308-82-9	32
2			2-Butenoic acid, 4-hydroxy-, met...	116	C5H8O3	004508-99-0	27
3			Ethyl hydrogen malonate	132	C5H8O4	001071-46-1	27
4			Octanoyl chloride	162	C8H15ClO	000111-64-8	14
5			Octanoic acid, silver(1+) salt	250	C8H15AgO2	024927-67-1	12



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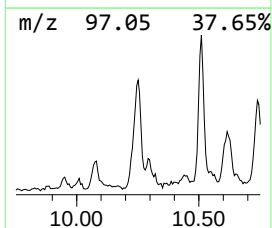
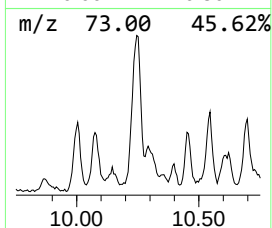
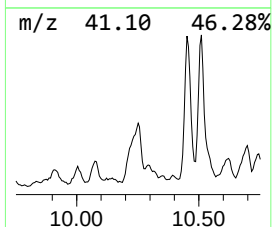
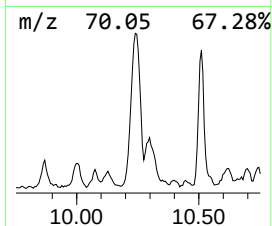
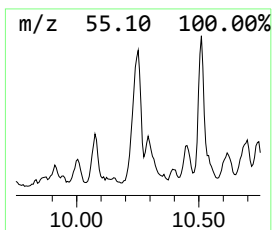
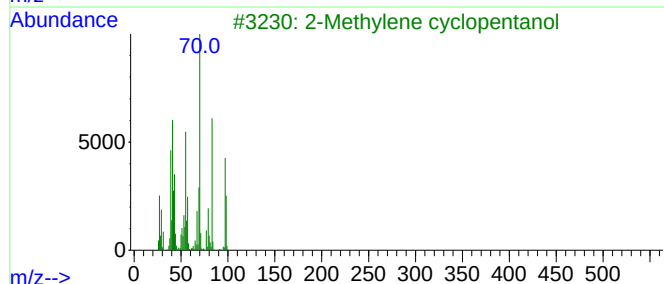
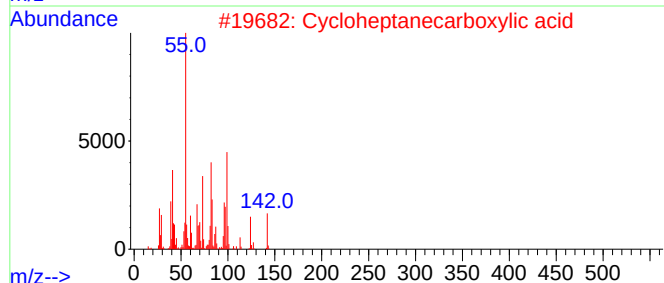
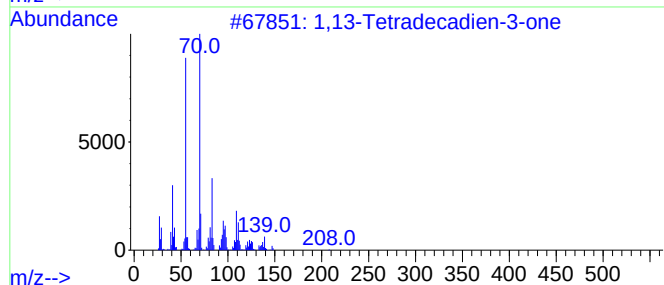
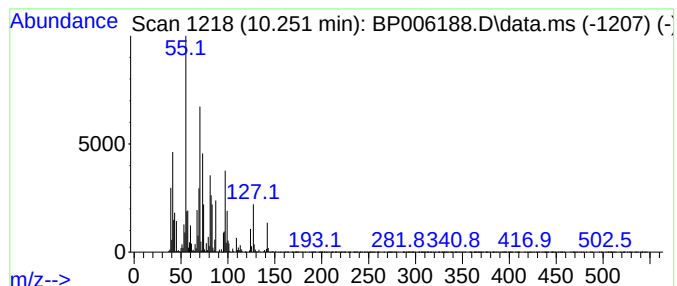
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 unknown10.251 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.251	12.88 ng	1045740	Naphthalene-d8	10.663

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,13-Tetradecadien-3-one	208	C14H24O	058879-40-6	35
2			Cycloheptanecarboxylic acid	142	C8H14O2	001460-16-8	25
3			2-Methylene cyclopentanol	98	C6H10O	020461-31-8	22
4			Acetamide, N-(3-methyl-2-buten-1...	127	C7H13NO	073286-69-8	14
5			1-Methylcyclohexylcarboxylic acid	142	C8H14O2	001123-25-7	14



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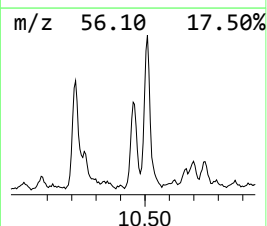
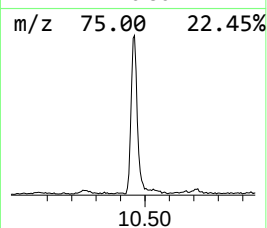
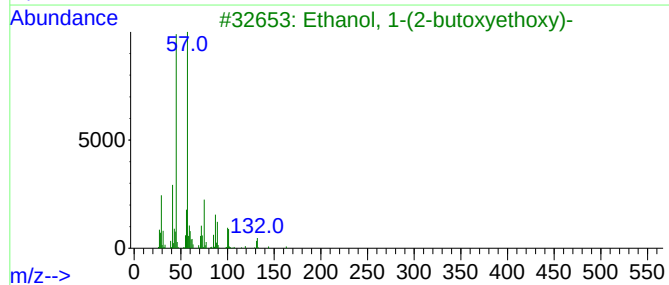
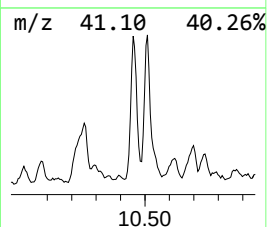
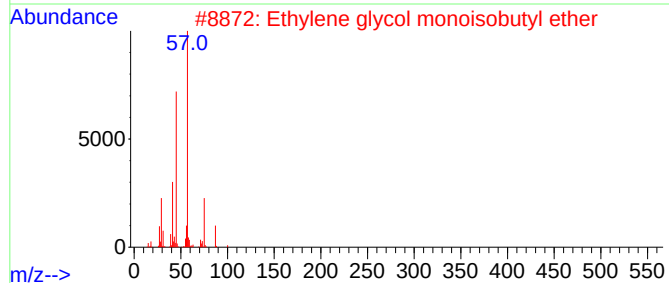
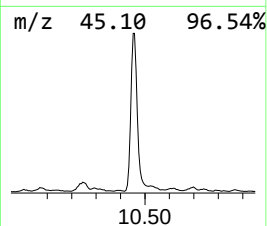
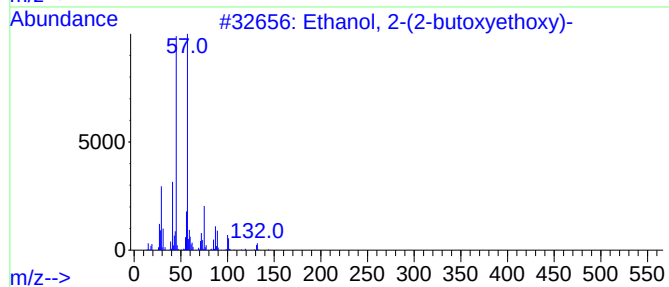
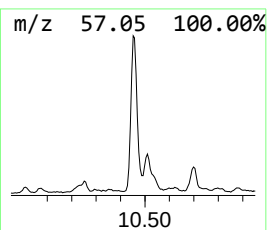
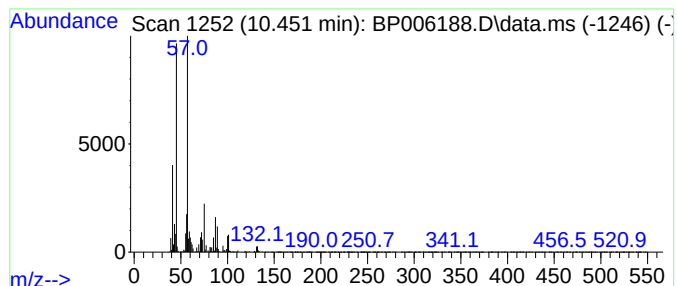
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Ethanol, 2-(2-butoxyethoxy)- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.451	13.24 ng	1074520	Naphthalene-d8	10.663

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
2			Ethylene glycol monoisobutyl ether	118	C6H14O2	004439-24-1	59
3			Ethanol, 1-(2-butoxyethoxy)-	162	C8H18O3	054446-78-5	50
4			3,3-Dimethylbutane-2-ol	102	C6H14O	000464-07-3	49
5			Methoxyacetic acid, 2-butyl ester	146	C7H14O3	070159-90-9	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071221\
Data File : BP006188.D
Acq On : 12 Jul 2021 19:04
Operator : CG/JU
Sample : M2997-01
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
RBR-200029

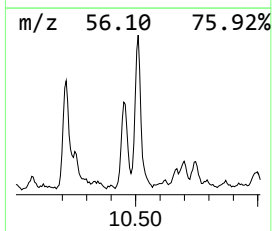
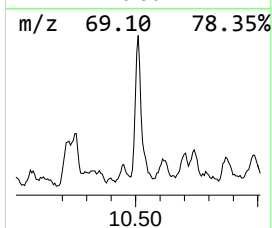
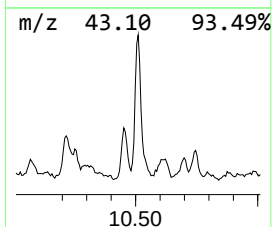
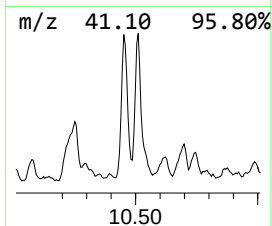
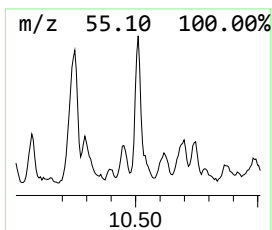
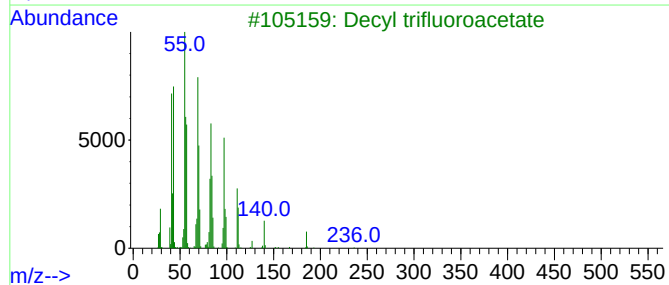
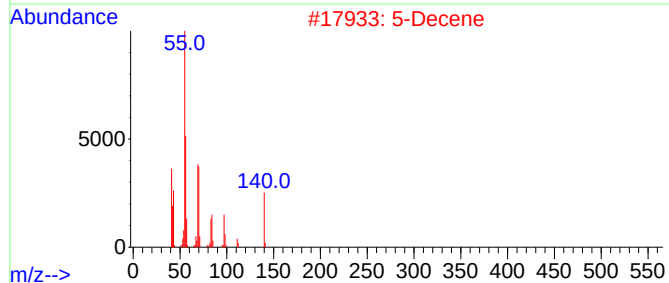
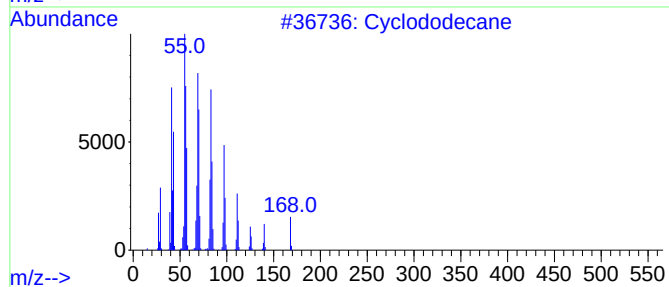
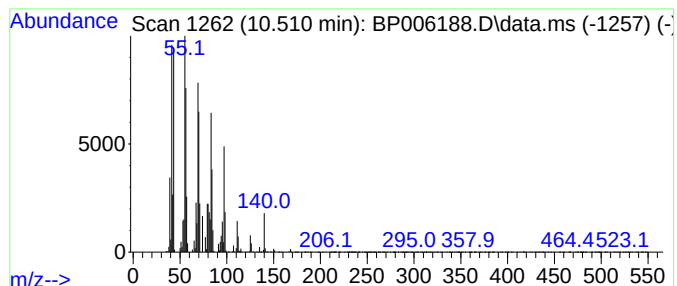
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Cyclododecane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.510	14.04 ng	1139740	Naphthalene-d8	10.663

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclododecane	168	C12H24	000294-62-2	91
2			5-Decene	140	C10H20	019689-19-1	90
3			Decyl trifluoroacetate	254	C12H21F3O2	000333-88-0	87
4			cis-3-Decene	140	C10H20	019398-86-8	76
5			1-Decene	140	C10H20	000872-05-9	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071221\
Data File : BP006188.D
Acq On : 12 Jul 2021 19:04
Operator : CG/JU
Sample : M2997-01
Misc :
ALS Vial : 7 Sample Multiplier: 1

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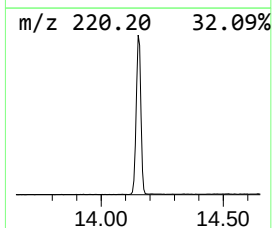
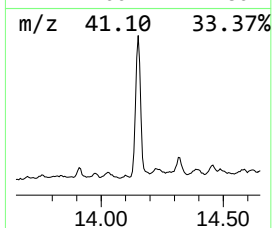
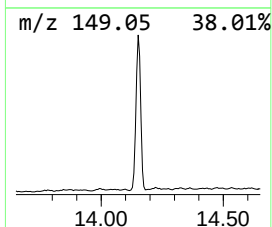
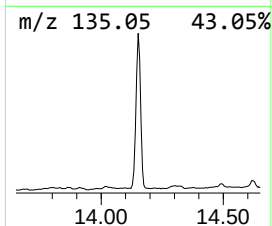
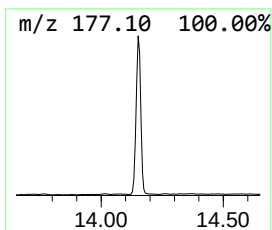
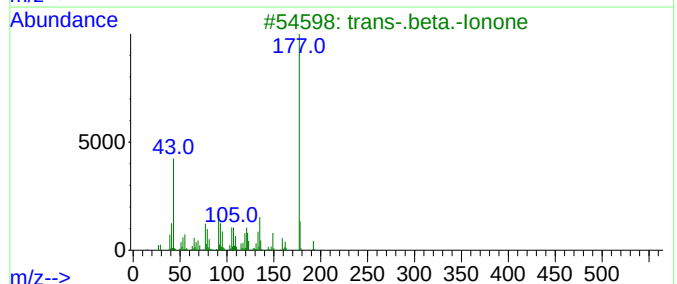
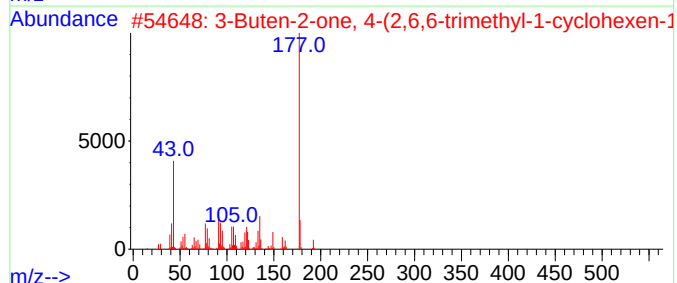
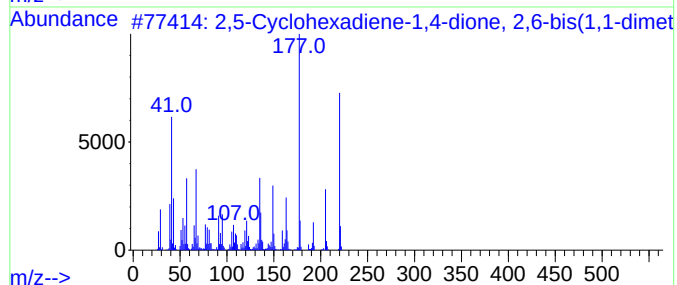
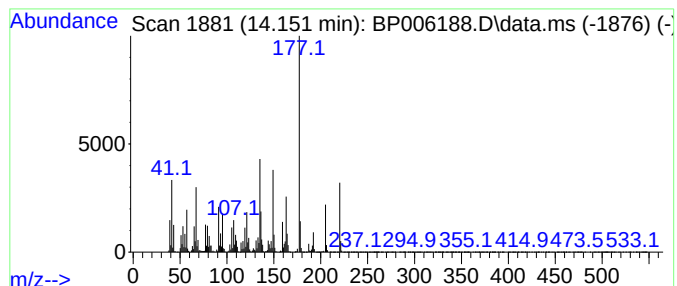
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 2,5-Cyclohexadiene-1,4-dione... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.151	78.99 ng	7808580	Acenaphthene-d10	14.498

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2,5-Cyclohexadiene-1,4-dione, 2,...	220	C14H20O2	000719-22-2	99
2			3-Buten-2-one, 4-(2,6,6-trimethy...	192	C13H20O	014901-07-6	50
3			trans-.beta.-Ionone	192	C13H20O	000079-77-6	50
4			Thiocoumarine, 4,4,6,7-tetramethyl-	220	C13H16OS	1000128-90-2	47
5			2H-1-Benzopyran, 3,5,6,8a-tetra...	192	C13H20O	041678-29-9	46



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071221\
Data File : BP006188.D
Acq On : 12 Jul 2021 19:04
Operator : CG/JU
Sample : M2997-01
Misc :
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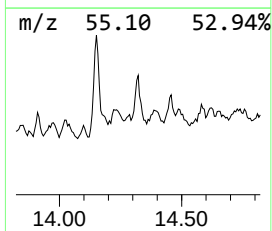
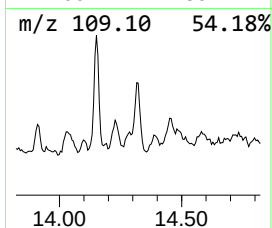
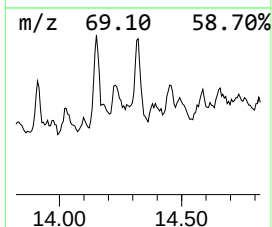
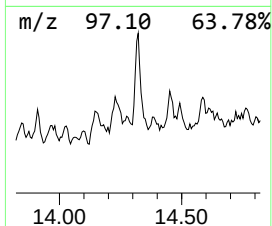
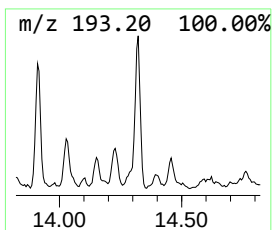
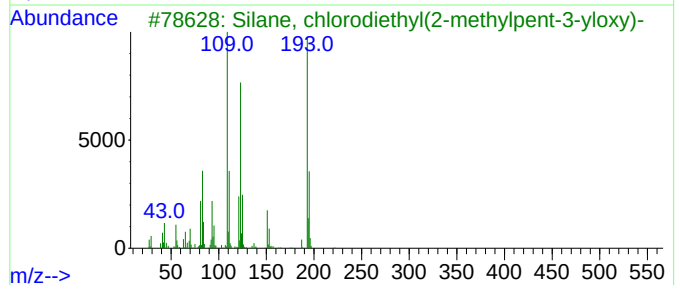
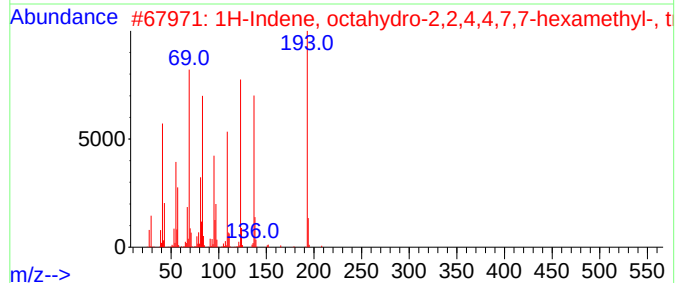
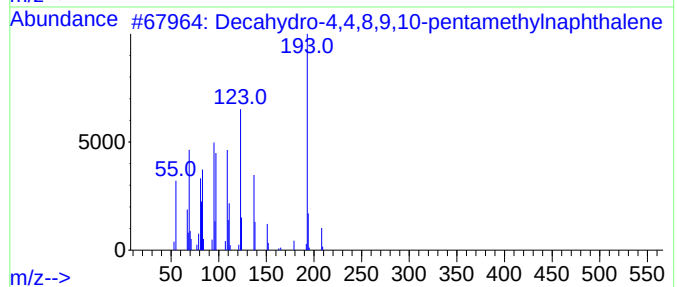
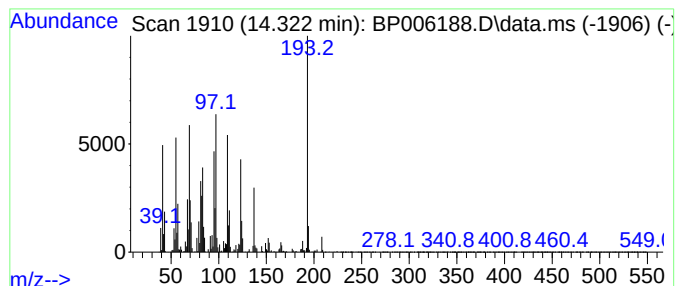
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Decahydro-4,4,8,9,10-pentam... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.322	8.13 ng	803966	Acenaphthene-d10	14.498

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Decahydro-4,4,8,9,10-pentamethyl...	208	C15H28	080655-44-3	64
2			1H-Indene, octahydro-2,2,4,4,7,7...	208	C15H28	054832-83-6	43
3			Silane, chlorodiethyl(2-methylpe...	222	C10H23ClOSi	1000363-79-1	40
4			5-Isobenzofurancarboxylic acid, ...	360	C21H28O5	022485-51-4	38
5			Silane, chlorodiethylhexyloxy-	222	C10H23ClOSi	1000363-74-4	25



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071221\
Data File : BP006188.D
Acq On : 12 Jul 2021 19:04
Operator : CG/JU
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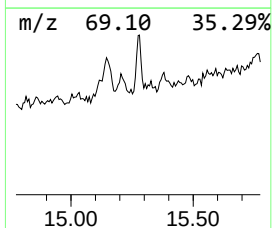
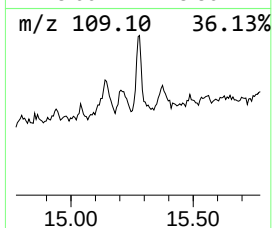
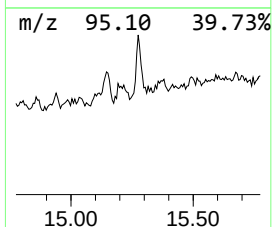
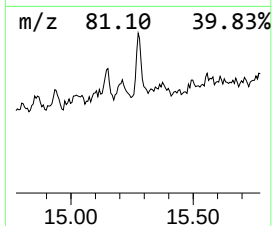
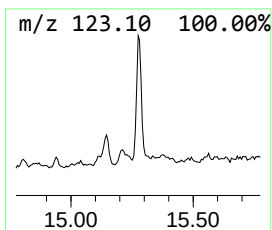
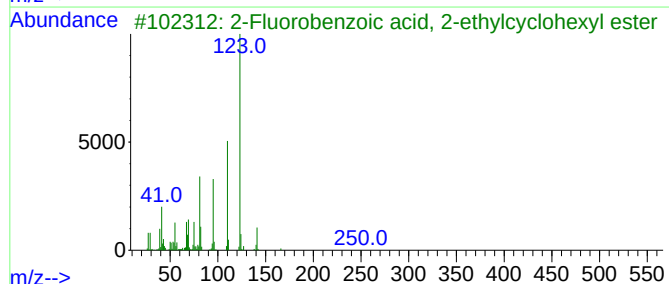
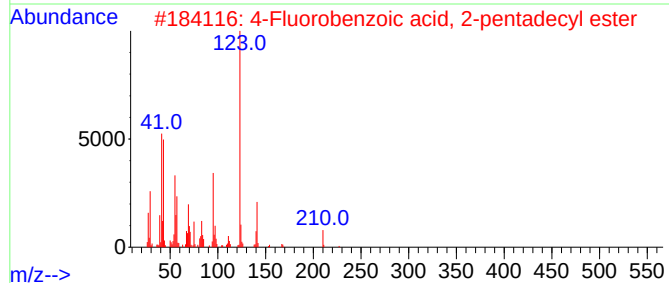
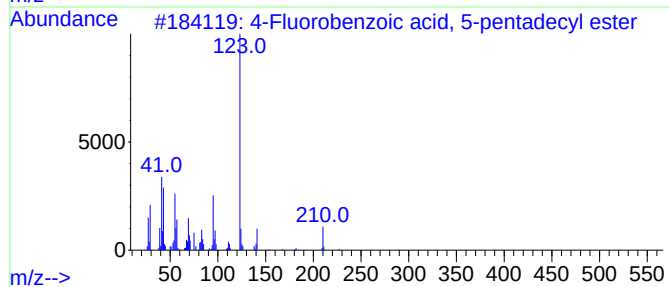
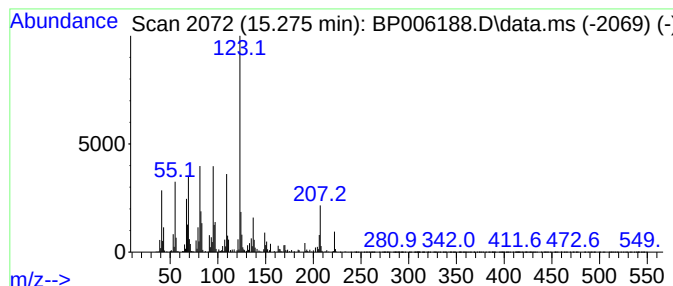
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 4-Fluorobenzoic acid, 5-pen... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.275	6.31 ng	624213	Acenaphthene-d10	14.498

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4-Fluorobenzoic acid, 5-pentadec...	350	C22H35F02	1000280-68-6	50
2			4-Fluorobenzoic acid, 2-pentadec...	350	C22H35F02	1000280-68-3	50
3			2-Fluorobenzoic acid, 2-ethylcyc...	250	C15H19F02	1000278-98-1	47
4			Decahydro-8a-ethyl-1,1,4a,6-tetr...	222	C16H30	1000100-23-6	46
5			(Phenylthio)acetic acid, cyclobu...	222	C12H14O2S	1000299-95-5	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071221\
Data File : BP006188.D
Acq On : 12 Jul 2021 19:04
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Sample : M2997-01
Misc :
ALS Vial : 7 Sample Multiplier: 1

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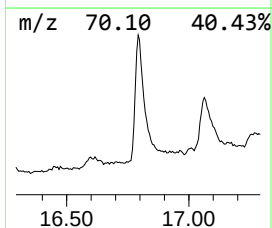
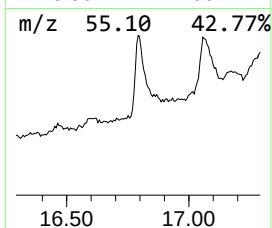
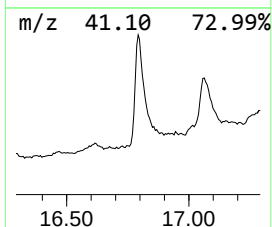
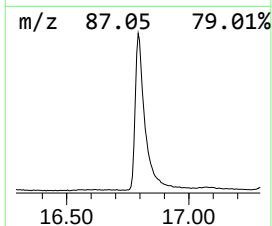
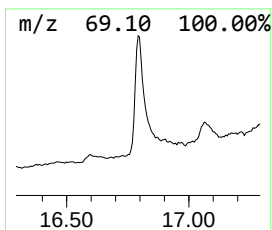
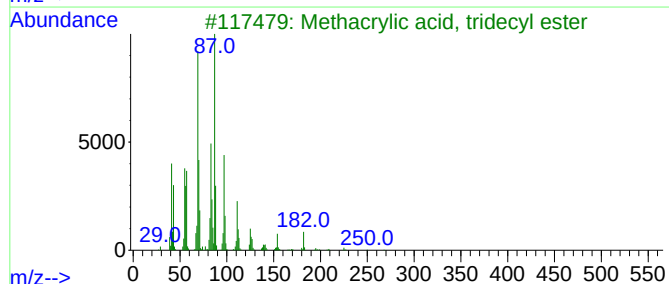
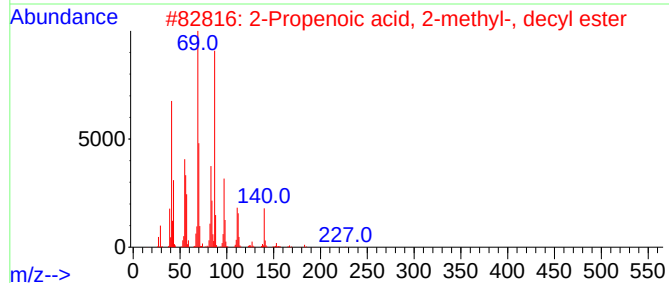
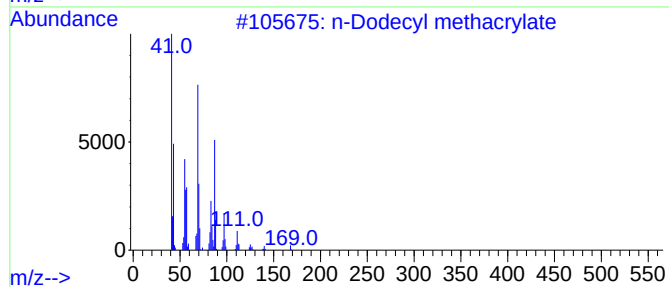
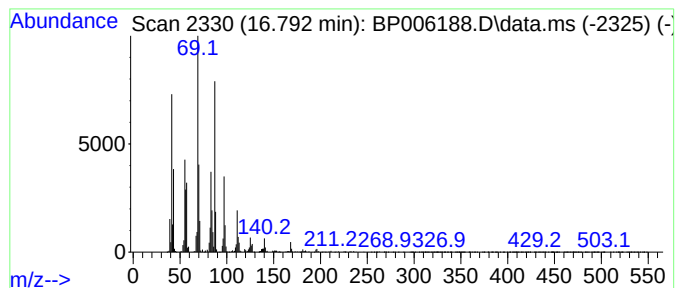
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 n-Dodecyl methacrylate Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.792	28.24 ng	7387570	Phenanthrene-d10	17.257

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Dodecyl methacrylate	254	C16H30O2	000142-90-5	91
2			2-Propenoic acid, 2-methyl-, dec...	226	C14H26O2	003179-47-3	87
3			Methacrylic acid, tridecyl ester	268	C17H32O2	1000340-28-9	62
4			Methacrylic acid, tetradecyl ester	282	C18H34O2	1000340-29-0	55
5			Cyclopropanecarboxylic acid, decy...	226	C14H26O2	054460-47-8	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071221\
Data File : BP006188.D
Acq On : 12 Jul 2021 19:04
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Misc :
ALS Vial : 7 Sample Multiplier: 1

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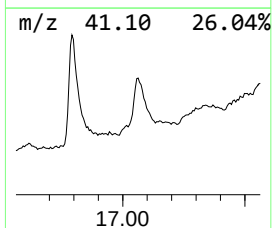
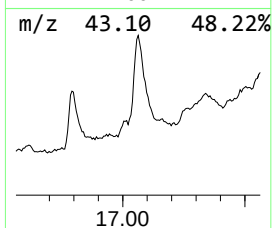
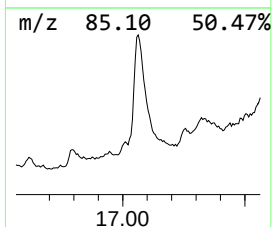
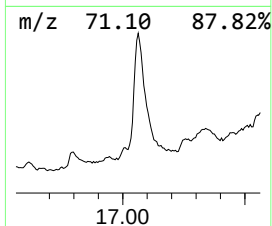
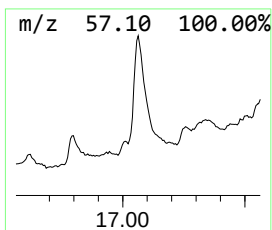
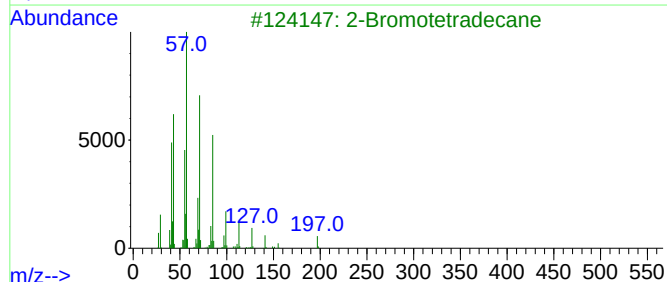
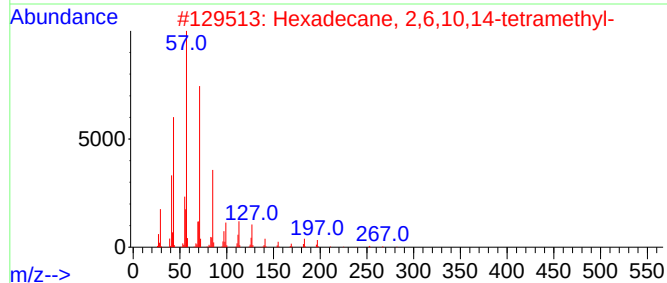
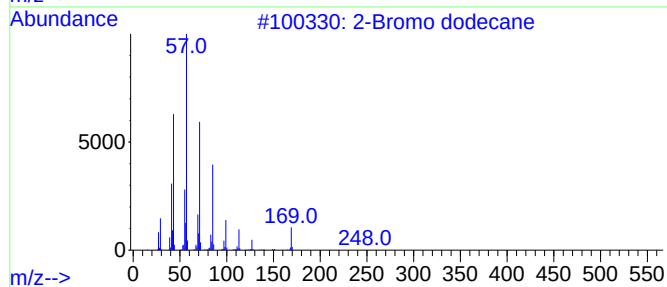
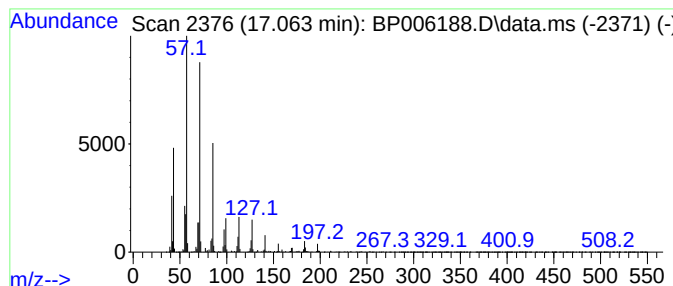
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 2-Bromo dodecane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.063	20.00 ng	5232570	Phenanthrene-d10	17.257

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		2-Bromo dodecane	248	C12H25Br	013187-99-0	91
2		Hexadecane, 2,6,10,14-tetramethyl-	282	C20H42	000638-36-8	91
3		2-Bromotetradecane	276	C14H29Br	074036-95-6	87
4		Heneicosane	296	C21H44	000629-94-7	86
5		Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071221\
Data File : BP006188.D
Acq On : 12 Jul 2021 19:04
Operator : CG/JU
Sample : M2997-01
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
RBR-200029

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
unknown7.722	7.722	8.2	ng	466695	1	7.863	1140220	20.0
1-Hexanol, 2-et...	7.940	189.7	ng	10812500	1	7.863	1140220	20.0
Hexanoic acid, ...	9.228	10.7	ng	610860	1	7.863	1140220	20.0
unknown9.351	9.351	7.7	ng	621223	2	10.663	1623200	20.0
unknown10.251	10.251	12.9	ng	1045740	2	10.663	1623200	20.0
Ethanol, 2-(2-b...	10.451	13.2	ng	1074520	2	10.663	1623200	20.0
Cyclododecane	10.510	14.0	ng	1139740	2	10.663	1623200	20.0
2,5-Cyclohexadi...	14.151	79.0	ng	7808580	3	14.498	1977090	20.0
Decahydro-4,4,8...	14.322	8.1	ng	803966	3	14.498	1977090	20.0
4-Fluorobenzoic...	15.275	6.3	ng	624213	3	14.498	1977090	20.0
n-Dodecyl metha...	16.792	28.2	ng	7387570	4	17.257	5232570	20.0
2-Bromo dodecane	17.063	20.0	ng	5232570	4	17.257	5232570	20.0