

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP012723\
 Data File : BP013625.D
 Acq On : 27 Jan 2023 19:24
 Operator : CG/JU
 Sample : 01320-01
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 TOTE-COMP-0126

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-BP012623.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BP013625.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.223	3	6	12	rVB	3384055	3947801	28.35%	6.058%
2	3.770	95	99	108	rBV	180264	232913	1.67%	0.357%
3	4.717	244	260	277	rBV4	56050	314532	2.26%	0.483%
4	5.217	339	345	355	rBV	527363	762035	5.47%	1.169%
5	5.687	418	425	436	rVB	2471144	3791061	27.22%	5.817%
6	7.117	662	668	676	rBV	235439	368772	2.65%	0.566%
7	7.228	682	687	692	rBV	111315	167369	1.20%	0.257%
8	7.293	692	698	708	rVB	1876733	3003953	21.57%	4.609%
9	8.152	837	844	850	rBV	703620	1093251	7.85%	1.677%
10	8.811	949	956	966	rVB	498733	760149	5.46%	1.166%
11	9.328	1035	1044	1054	rBV	2370064	3865039	27.76%	5.931%
12	10.981	1316	1325	1332	rBV	879666	1478614	10.62%	2.269%
13	11.634	1429	1436	1447	rBV	248903	447521	3.21%	0.687%
14	12.581	1588	1597	1615	rBV	440692	959272	6.89%	1.472%
15	13.416	1730	1739	1751	rBV	6513705	9896703	71.07%	15.186%
16	14.722	1956	1961	1967	rBV	362329	525369	3.77%	0.806%
17	14.787	1967	1972	1980	rVB	1300233	1983148	14.24%	3.043%
18	16.145	2198	2203	2212	rVB2	332266	494344	3.55%	0.759%
19	16.275	2218	2225	2238	rBV	4298381	6393760	45.91%	9.811%
20	16.540	2264	2270	2276	rBV	110985	157298	1.13%	0.241%
21	17.010	2345	2350	2361	rBV	320904	581843	4.18%	0.893%
22	17.292	2393	2398	2412	rVB	125065	214138	1.54%	0.329%
23	17.545	2435	2441	2449	rBV	1639032	2385566	17.13%	3.660%
24	18.398	2581	2586	2597	rBV	422273	677632	4.87%	1.040%
25	19.622	2790	2794	2803	rBV	194383	295968	2.13%	0.454%
26	20.081	2866	2872	2884	rVB	10547483	13925425	100.00%	21.367%
27	21.298	3075	3079	3086	rBV	427368	613063	4.40%	0.941%
28	21.622	3129	3134	3140	rBV	2110784	2887569	20.74%	4.431%
29	24.216	3568	3575	3584	rVB2	1405199	2947666	21.17%	4.523%

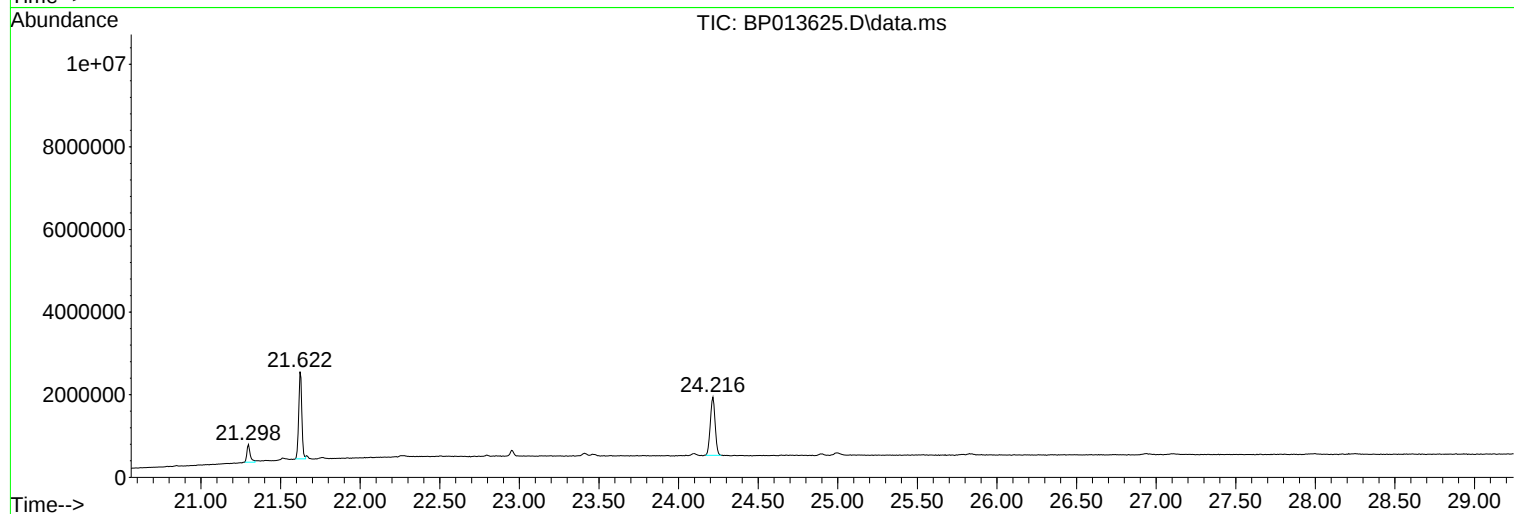
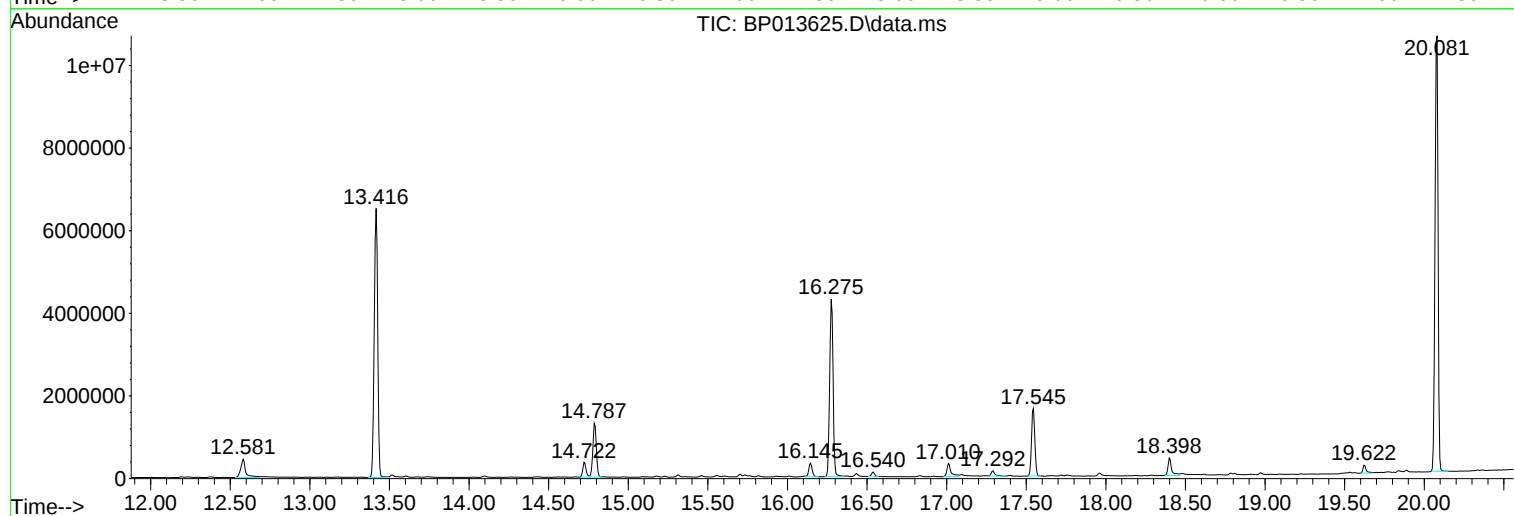
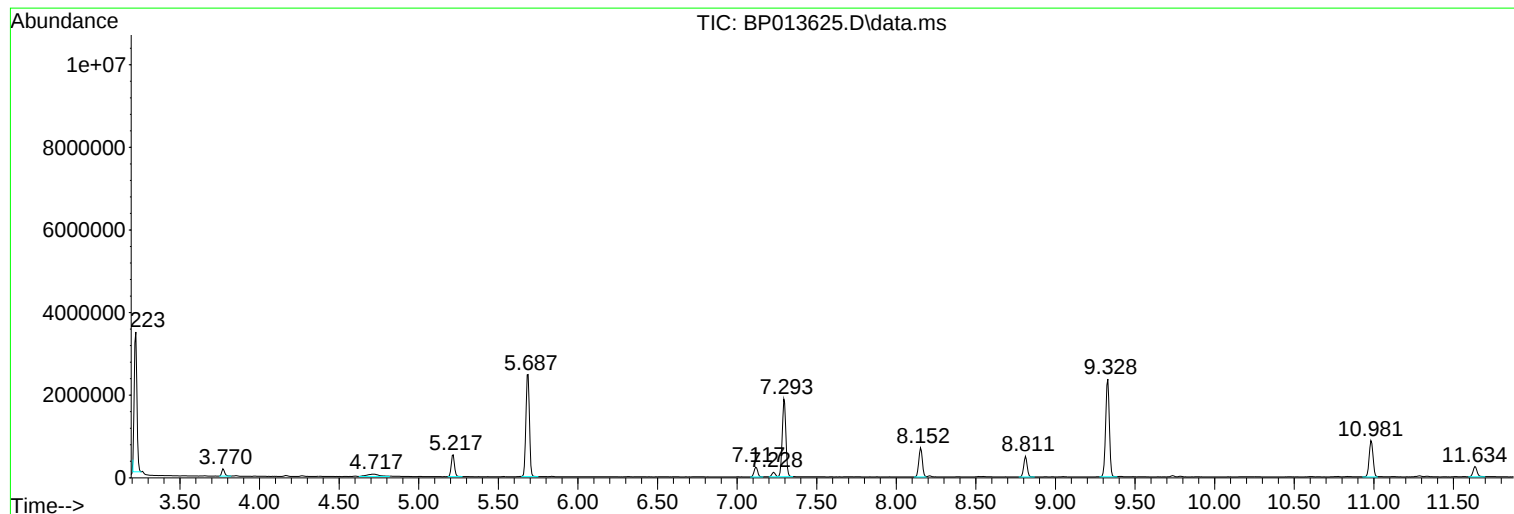
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP012623.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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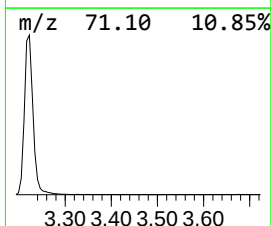
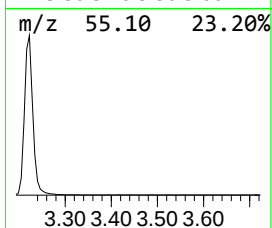
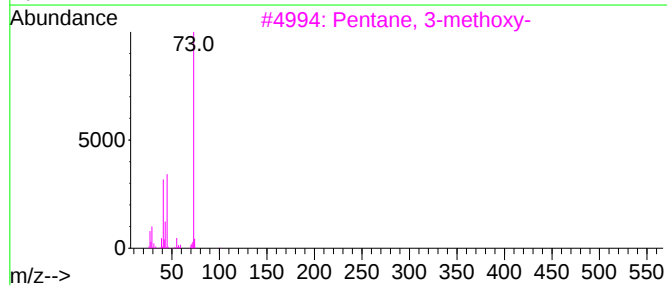
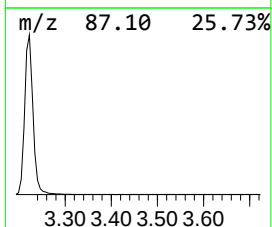
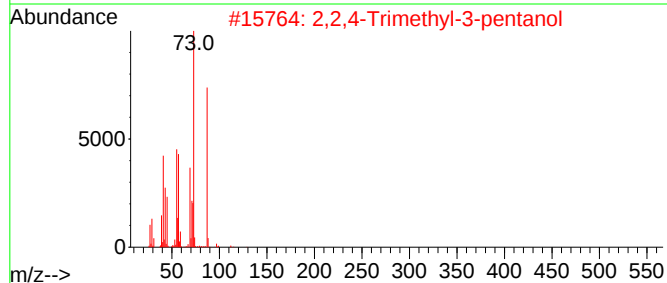
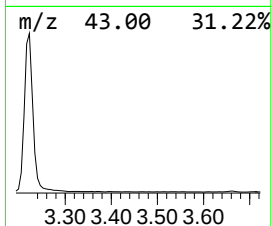
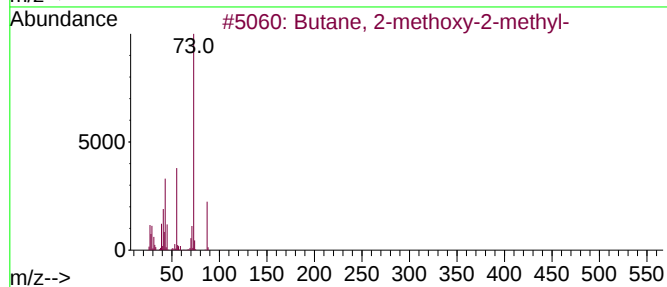
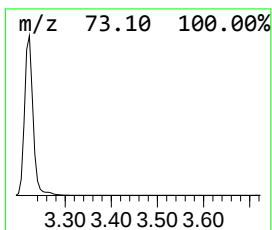
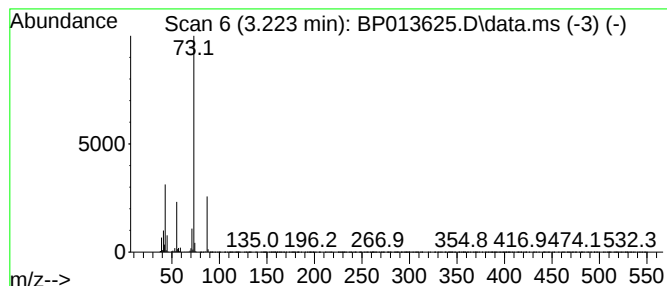
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.223	72.22 ng	3947800	1,4-Dichlorobenzene-d4	8.152

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2			2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39
3			Pentane, 3-methoxy-	102	C6H14O	036839-67-5	23
4			1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	23
5			Silane, tetramethyl-	88	C4H12Si	000075-76-3	17



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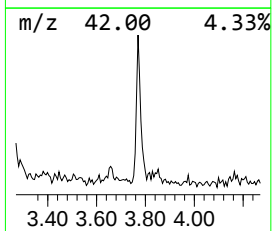
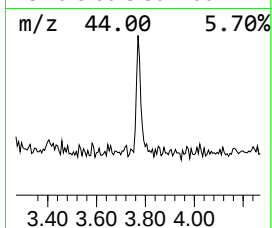
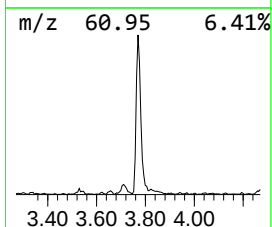
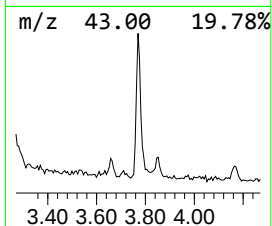
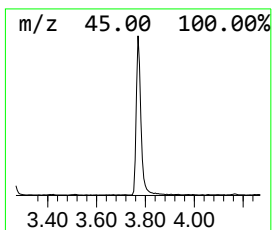
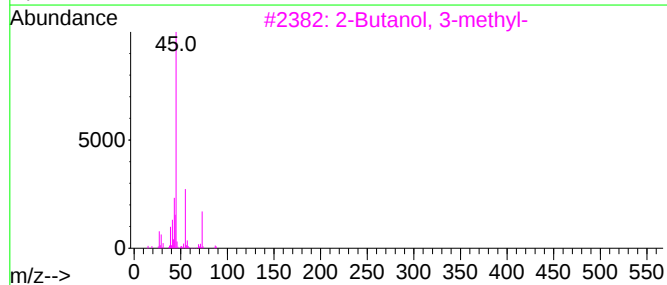
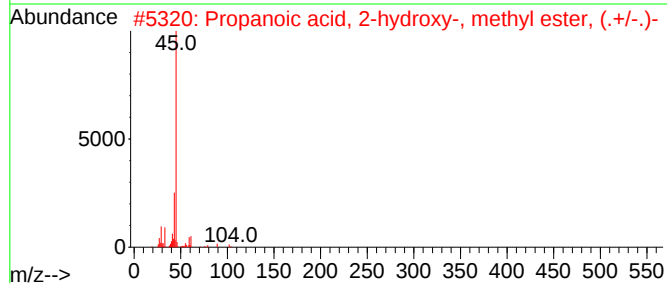
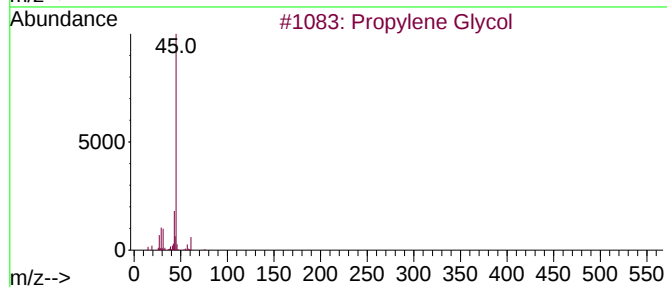
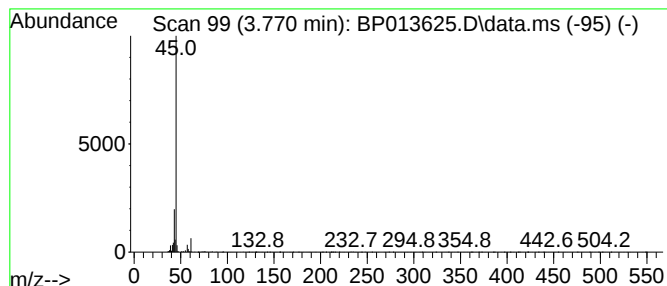
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP012623.M
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Propylene Glycol Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.770	4.26 ng	232913	1,4-Dichlorobenzene-d4	8.152

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Propylene Glycol	76	C3H8O2	000057-55-6	90
2			Propanoic acid, 2-hydroxy-, meth...	104	C4H8O3	000547-64-8	40
3			2-Butanol, 3-methyl-	88	C5H12O	000598-75-4	9
4			Propane, 2-ethoxy-	88	C5H12O	000625-54-7	9
5			R-(-)-1,2-propanediol	76	C3H8O2	004254-14-2	9



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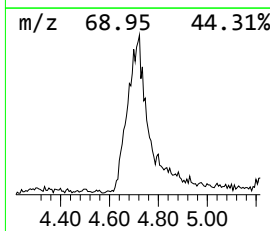
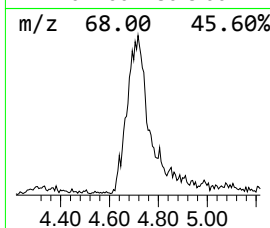
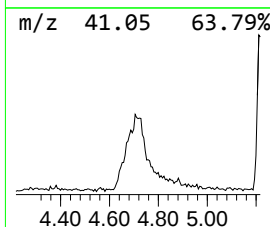
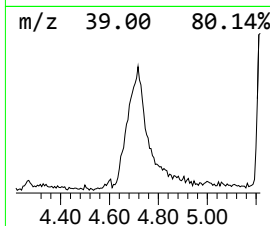
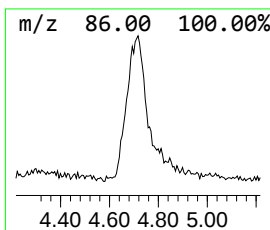
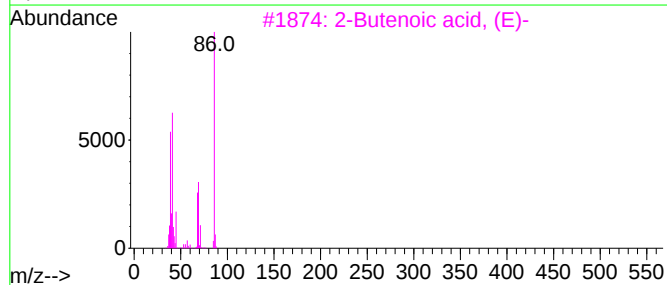
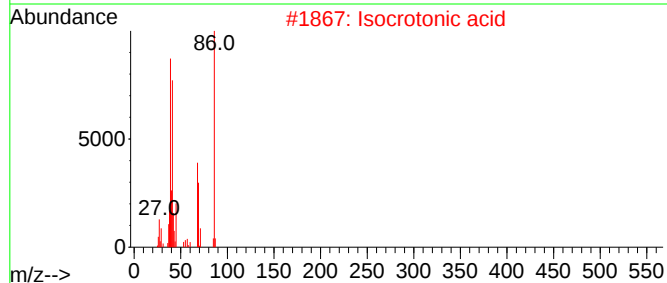
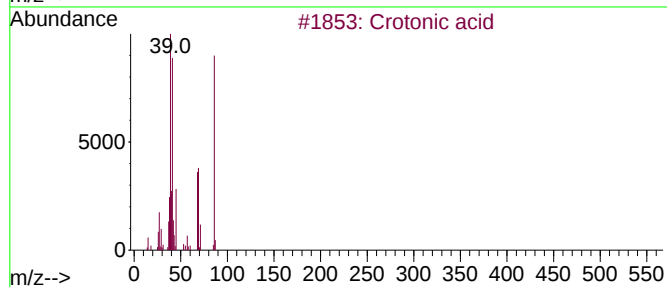
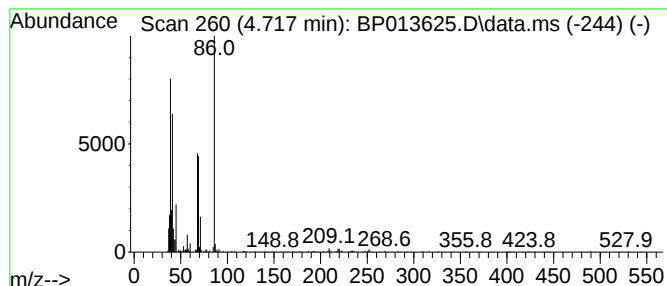
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Crotonic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.717	5.75 ng	314532	1,4-Dichlorobenzene-d4	8.152

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Crotonic acid	86	C4H6O2	003724-65-0	94
2			Isocrotonic acid	86	C4H6O2	000503-64-0	91
3			2-Butenoic acid, (E)-	86	C4H6O2	000107-93-7	90
4			2-Propenoic acid, 2-methyl-	86	C4H6O2	000079-41-4	87
5			3-Butenoic acid	86	C4H6O2	000625-38-7	72



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ALS Vial : 8 Sample Multiplier: 1

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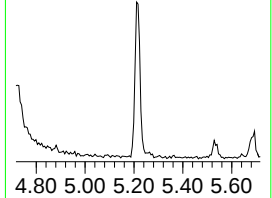
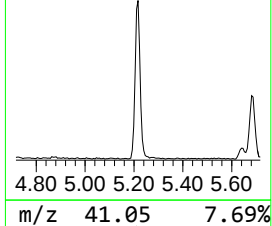
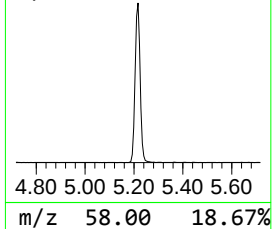
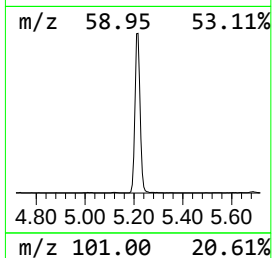
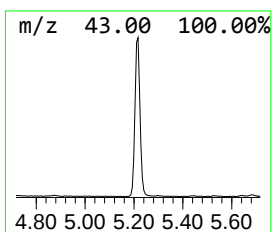
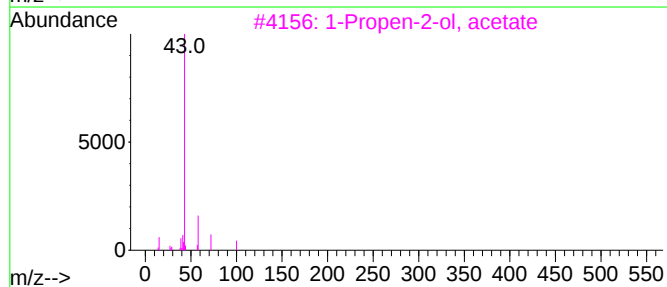
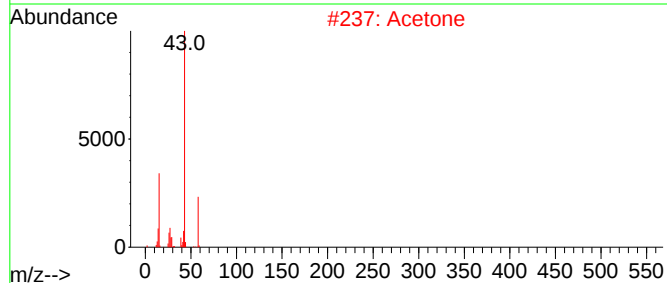
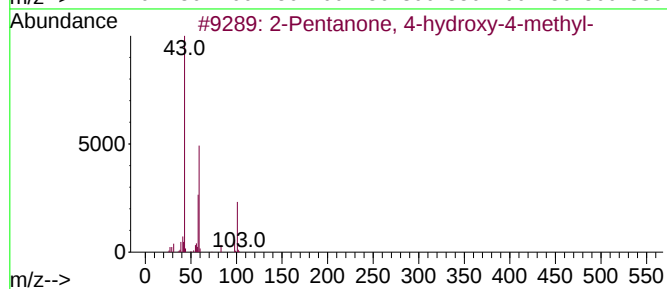
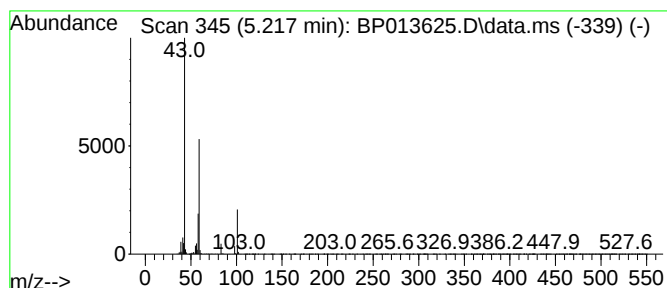
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Peak Number 4 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.217	13.94 ng	762035	1,4-Dichlorobenzene-d4	8.152

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	53
2			Acetone	58	C3H6O	000067-64-1	10
3			1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10
4			Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9
5			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	9



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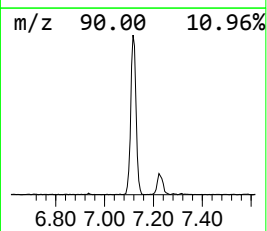
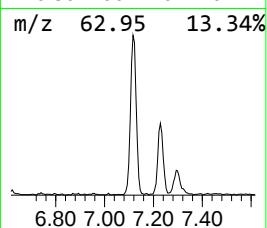
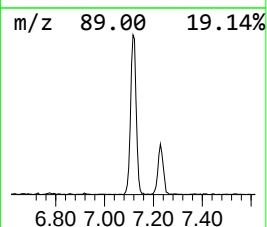
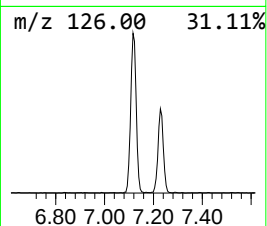
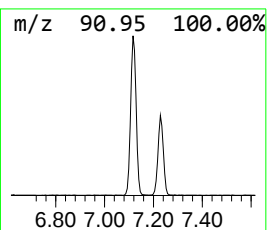
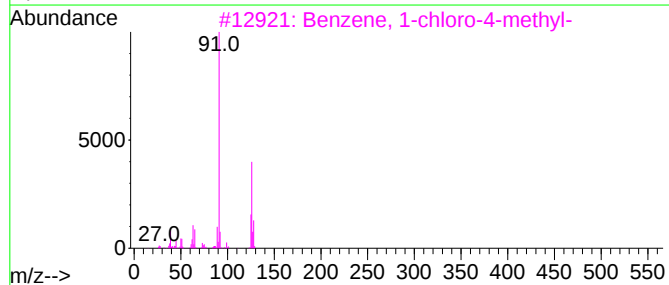
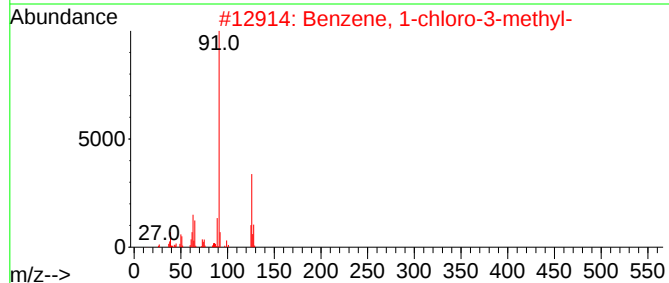
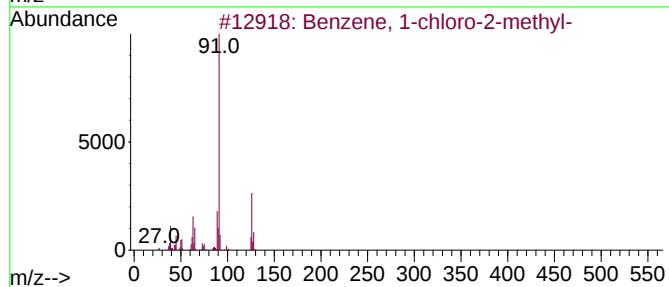
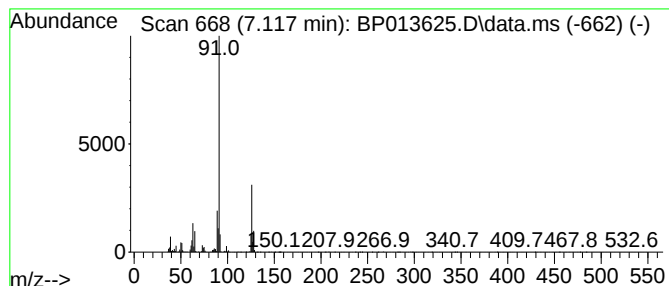
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Peak Number 5 Benzene, 1-chloro-2-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.117	6.75 ng	368772	1,4-Dichlorobenzene-d4	8.152

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-chloro-2-methyl-	126	C7H7Cl	000095-49-8	96
2			Benzene, 1-chloro-3-methyl-	126	C7H7Cl	000108-41-8	95
3			Benzene, 1-chloro-4-methyl-	126	C7H7Cl	000106-43-4	81
4			Benzyl chloride	126	C7H7Cl	000100-44-7	74
5			Benzaldehyde, 4-methyl-	120	C8H8O	000104-87-0	32



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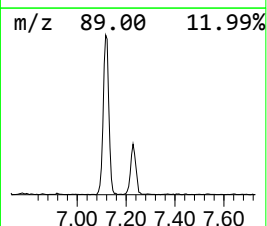
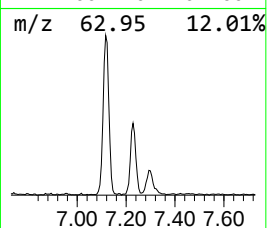
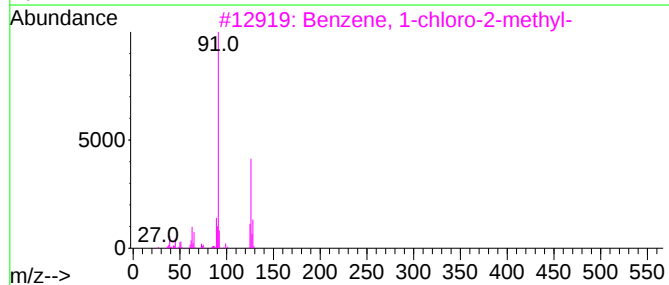
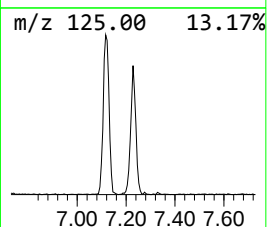
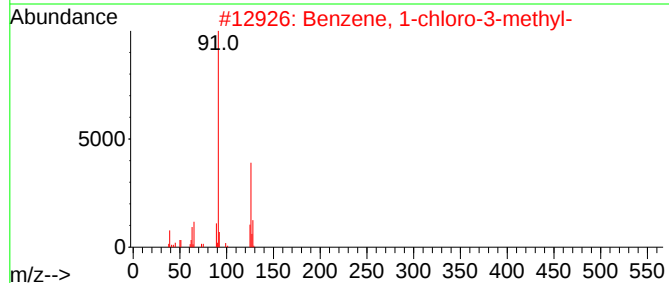
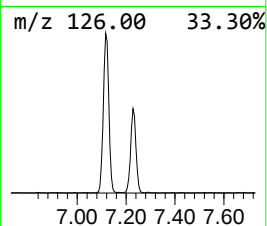
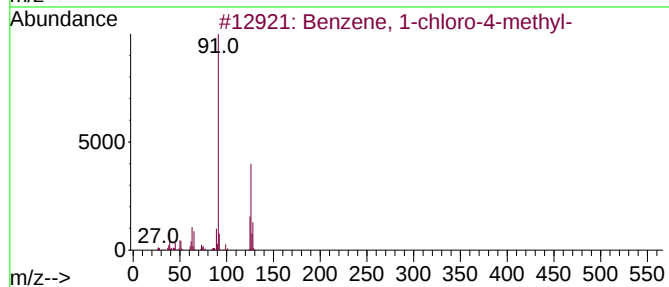
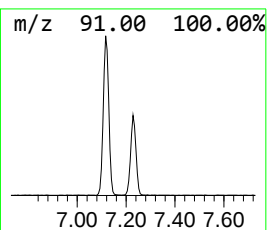
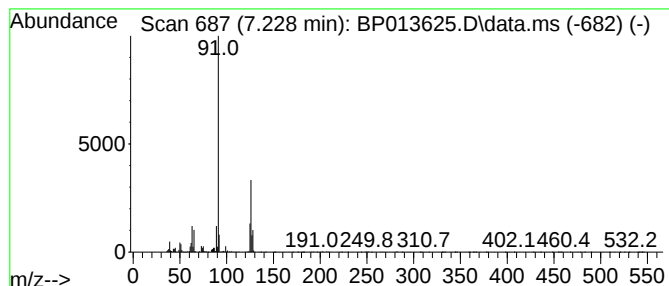
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Benzene, 1-chloro-4-methyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.228	3.06 ng	167369	1,4-Dichlorobenzene-d4	8.152

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-chloro-4-methyl-	126	C7H7Cl	000106-43-4	97
2			Benzene, 1-chloro-3-methyl-	126	C7H7Cl	000108-41-8	96
3			Benzene, 1-chloro-2-methyl-	126	C7H7Cl	000095-49-8	94
4			Benzyl chloride	126	C7H7Cl	000100-44-7	90
5			Benzene, (bromomethyl)-	170	C7H7Br	000100-39-0	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP012723\
Data File : BP013625.D
Acq On : 27 Jan 2023 19:24
Operator : CG/JU
Sample : 01320-01
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
TOTE-COMP-0126

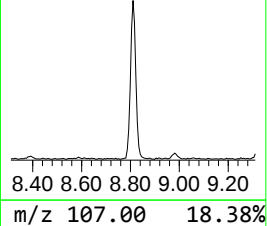
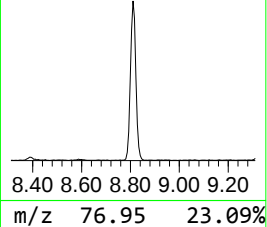
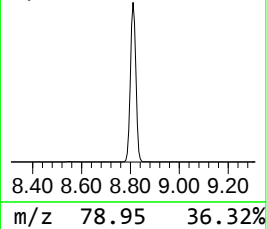
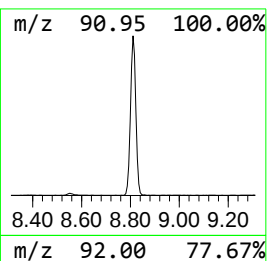
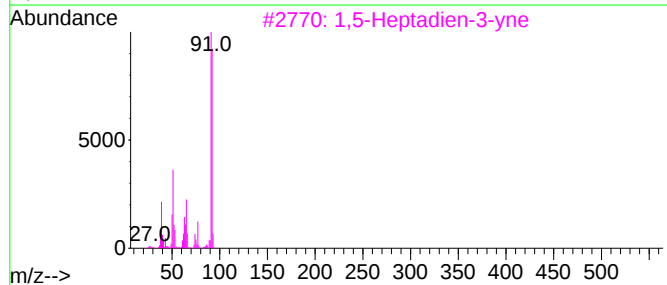
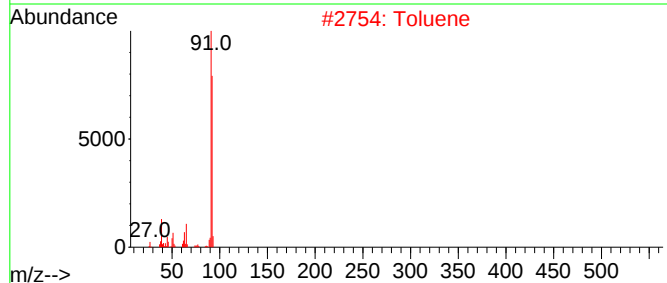
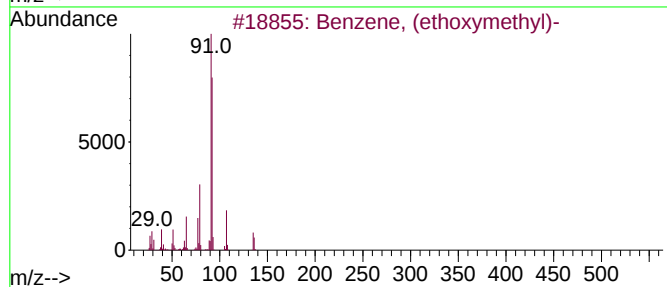
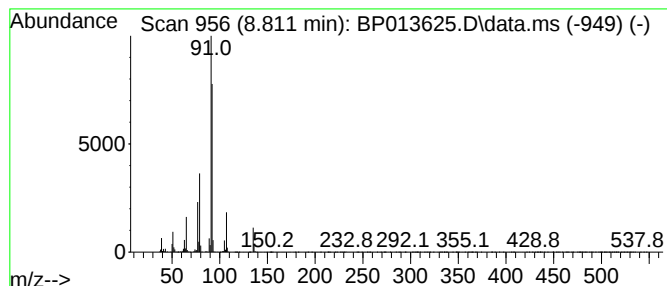
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Benzene, (ethoxymethyl)- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.811	13.91 ng	760149	1,4-Dichlorobenzene-d4	8.152

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, (ethoxymethyl)-	136	C9H12O	000539-30-0	91
2			Toluene	92	C7H8	000108-88-3	58
3			1,5-Heptadien-3-yne	92	C7H8	003511-27-1	58
4			Benzene, 1,1'-[oxybis(methylene)...]	198	C14H14O	000103-50-4	53
5			Benzeneacetamide	135	C8H9NO	000103-81-1	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP012723\
Data File : BP013625.D
Acq On : 27 Jan 2023 19:24
Operator : CG/JU
Sample : 01320-01
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
TOTE-COMP-0126

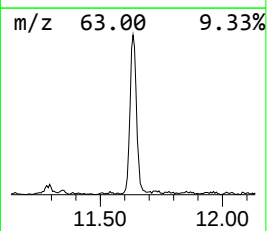
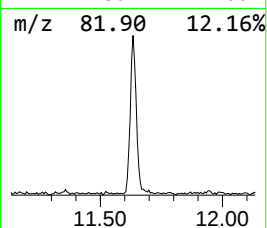
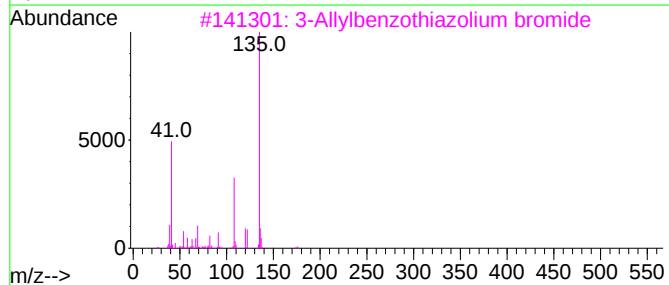
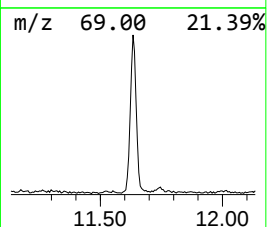
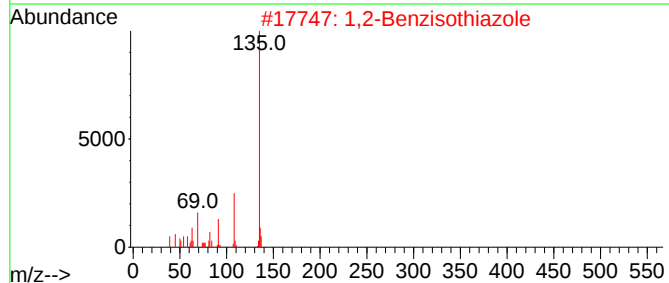
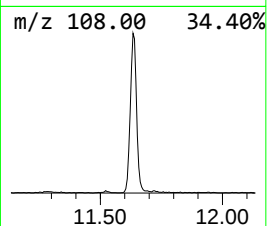
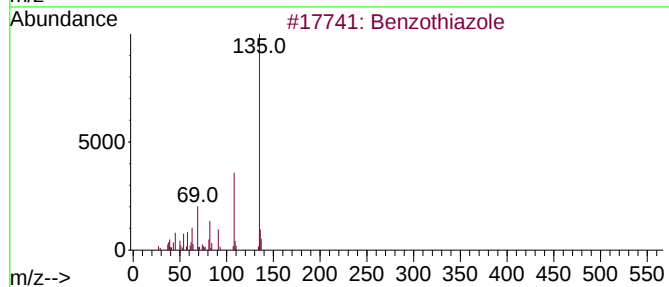
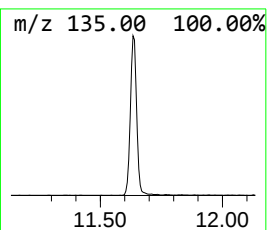
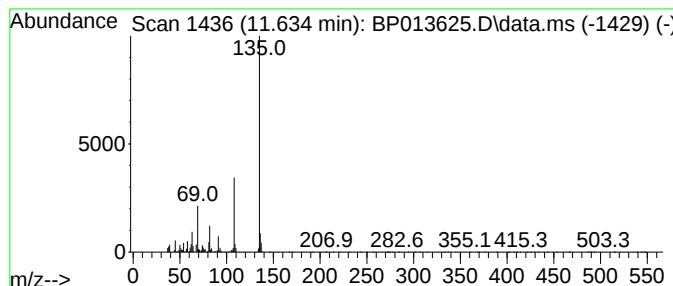
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Benzothiazole Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.634	6.05 ng	447521	Naphthalene-d8	10.981

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzothiazole	135	C7H5NS	000095-16-9	95
2			1,2-Benzisothiazole	135	C7H5NS	000272-16-2	90
3			3-Allylbenzothiazolium bromide	255	C10H10BrNS	016407-55-9	72
4			Thieno[2,3-c]pyridine	135	C7H5NS	000272-12-8	72
5			Adenine	135	C5H5N5	000073-24-5	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP012723\
Data File : BP013625.D
Acq On : 27 Jan 2023 19:24
Operator : CG/JU
Sample : 01320-01
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
TOTE-COMP-0126

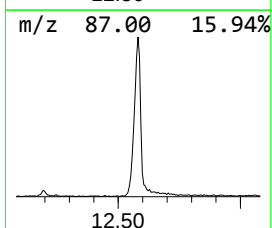
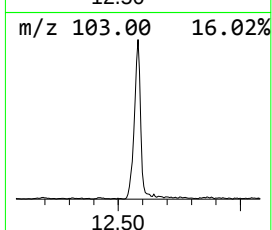
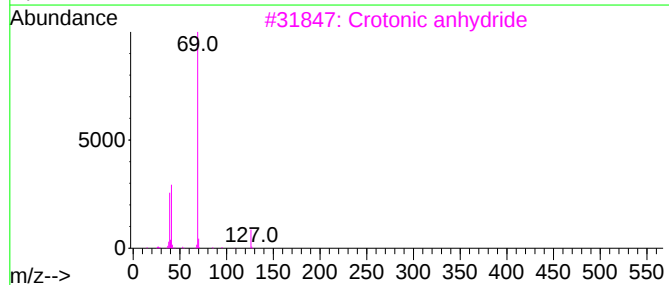
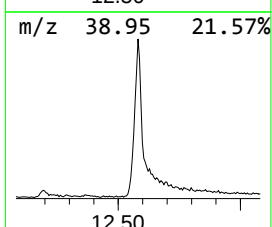
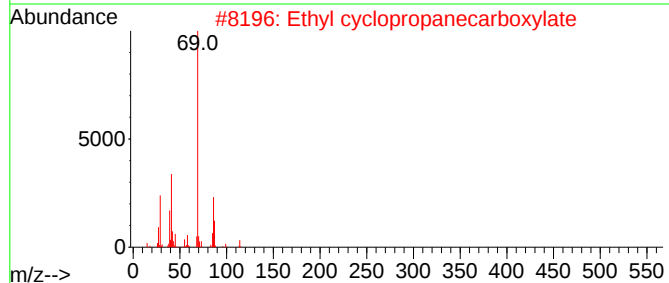
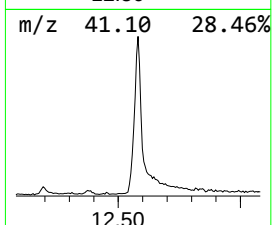
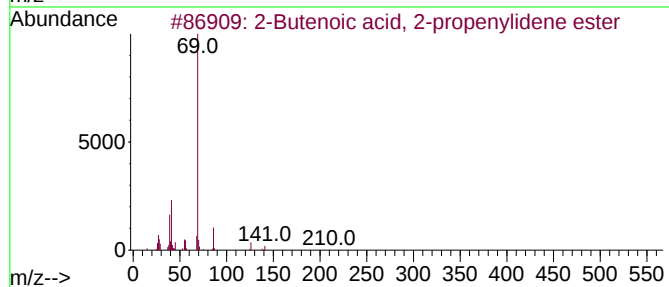
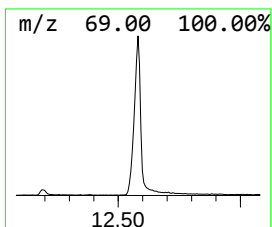
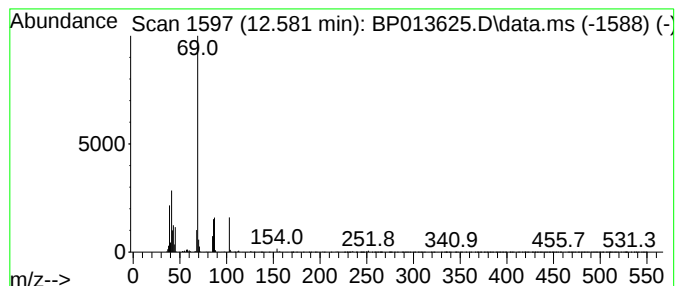
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 unknown12.581 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.581	12.98 ng	959272	Naphthalene-d8	10.981

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Butenoic acid, 2-propenylidene...	210	C11H14O4	055030-70-1	42
2			Ethyl cyclopropanecarboxylate	114	C6H10O2	004606-07-9	39
3			Crotonic anhydride	154	C8H10O3	000623-68-7	37
4			Isoamyl crotonate	156	C9H16O2	1010429-17-3	33
5			Oxazole	69	C3H3NO	000288-42-6	33



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP012723\
Data File : BP013625.D
Acq On : 27 Jan 2023 19:24
Operator : CG/JU
Sample : 01320-01
Misc :
ALS Vial : 8 Sample Multiplier: 1

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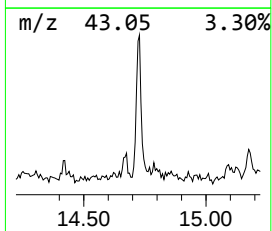
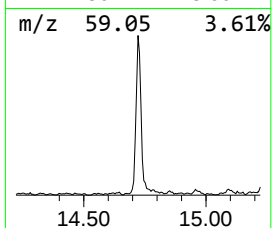
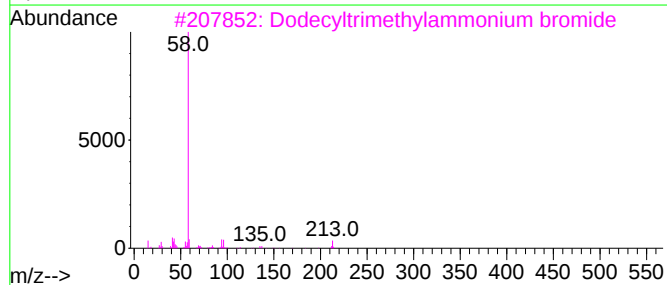
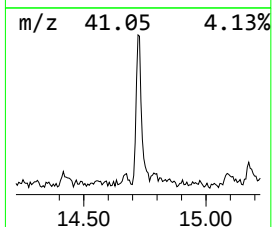
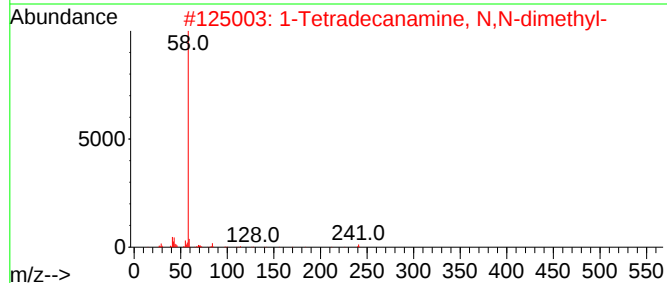
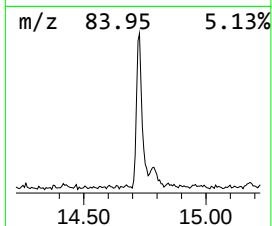
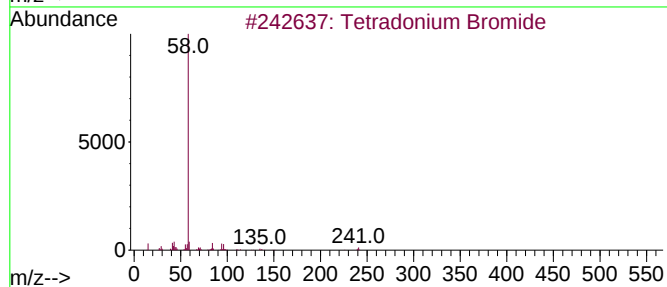
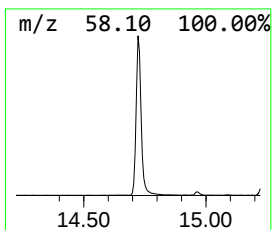
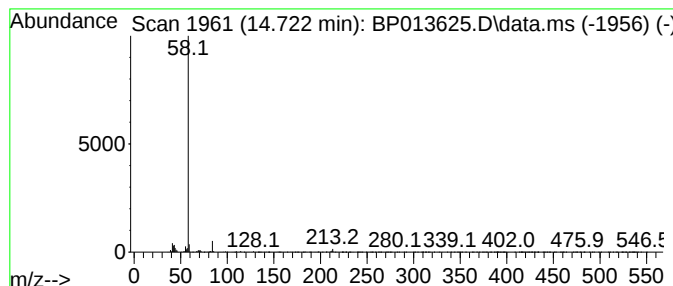
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Tetradonium Bromide Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.722	5.30 ng	525369	Acenaphthene-d10	14.787

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetradonium Bromide	335	C17H38BrN	001119-97-7	72
2			1-Tetradecanamine, N,N-dimethyl-	241	C16H35N	000112-75-4	72
3			Dodecyltrimethylammonium bromide	307	C15H34BrN	001119-94-4	72
4			Dimethyl palmitamine	269	C18H39N	000112-69-6	72
5			1-Decanaminium, N,N,N-trimethyl-...	235	C13H30ClN	010108-87-9	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP012723\
Data File : BP013625.D
Acq On : 27 Jan 2023 19:24
Operator : CG/JU
Sample : 01320-01
Misc :
ALS Vial : 8 Sample Multiplier: 1

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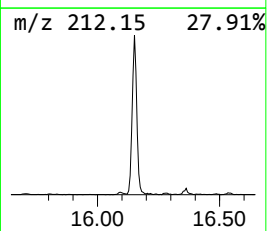
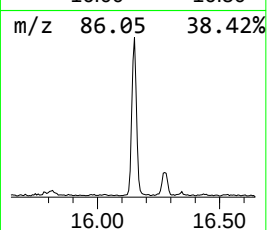
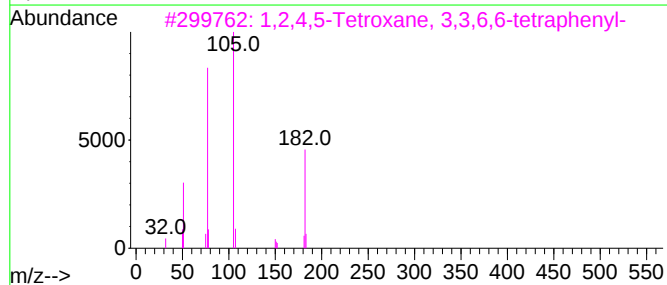
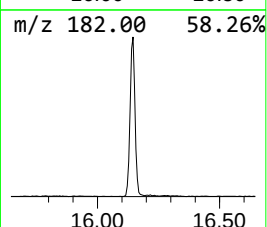
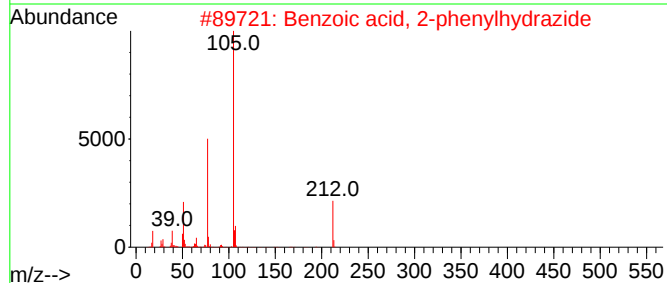
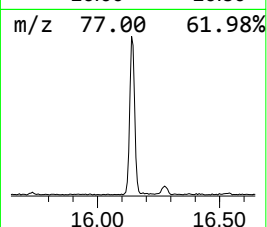
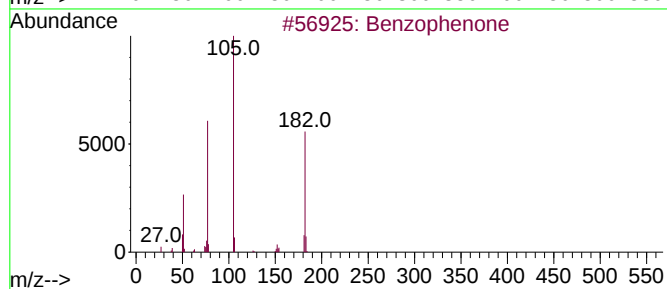
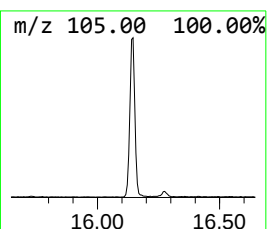
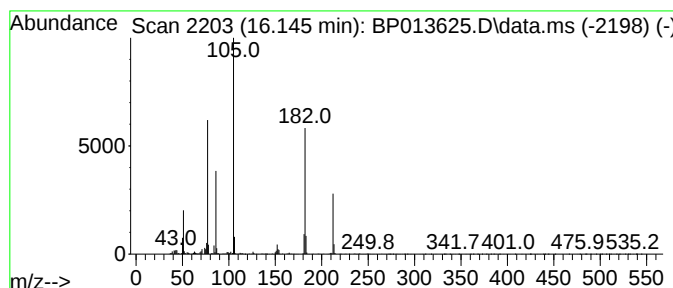
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Benzophenone Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.145	4.99 ng	494344	Acenaphthene-d10	14.787

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	96
2			Benzoic acid, 2-phenylhydrazide	212	C13H12N2O	000532-96-7	43
3			1,2,4,5-Tetroxane, 3,3,6,6-tetra...	396	C26H20O4	016204-36-7	40
4			Benzoic acid, 1,3-dioxan-5-yl ester	208	C11H12O4	049784-60-3	38
5			Phenyl 2-(2'-furanyl)cyclopropyl...	212	C14H12O2	1000140-13-9	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP012723\
Data File : BP013625.D
Acq On : 27 Jan 2023 19:24
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Misc :
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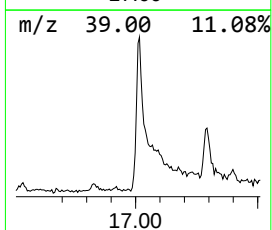
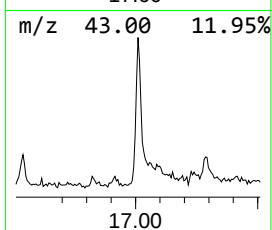
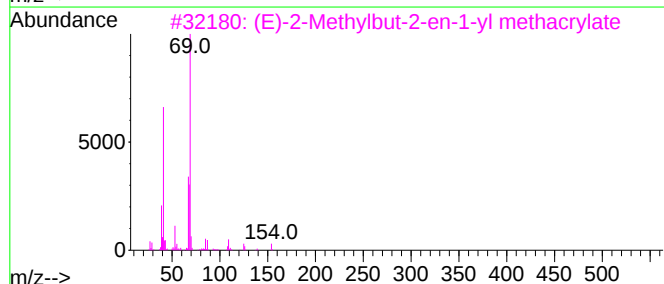
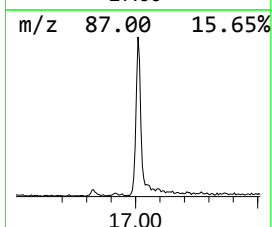
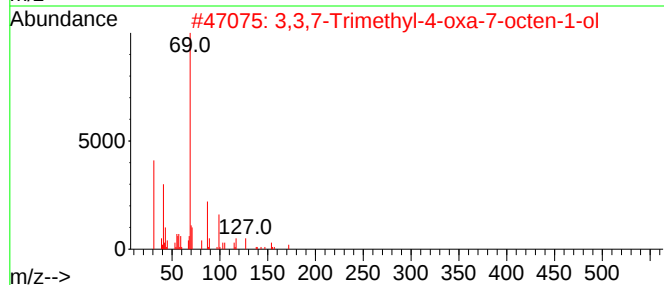
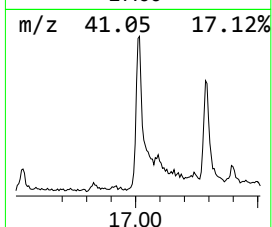
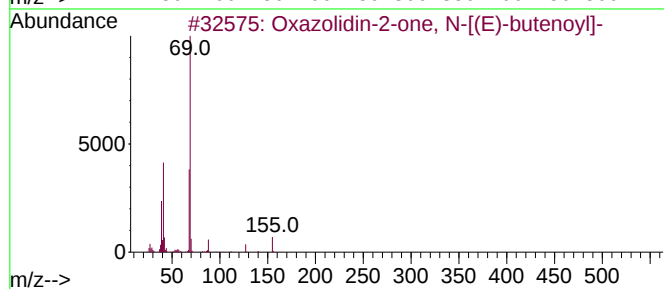
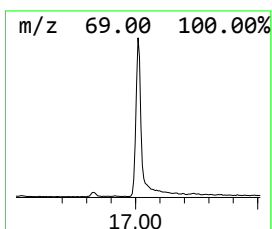
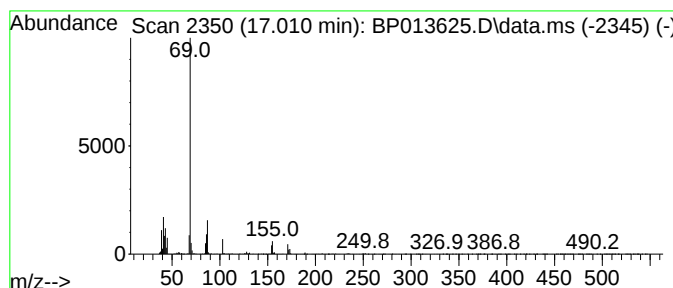
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Oxazolidin-2-one, N-[(E)-bu... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.	
17.010	4.88 ng	581843	Phenanthrene-d10	17.545	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Oxazolidin-2-one, N-[(E)-butenoyl]-	155	C7H9NO3	1010156-45-5	52
2	3,3,7-Trimethyl-4-oxa-7-octen-1-ol	172	C10H20O2	069161-47-3	38
3	(E)-2-Methylbut-2-en-1-yl methac...	154	C9H14O2	088142-95-4	9
4	4-Undecanone, 7-ethyl-2-methyl-	212	C14H28O	006976-00-7	9
5	Ethyl cyclopropanecarboxylate	114	C6H10O2	004606-07-9	9



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP012723\
Data File : BP013625.D
Acq On : 27 Jan 2023 19:24
Operator : CG/JU
Sample : 01320-01
Misc :
ALS Vial : 8 Sample Multiplier: 1

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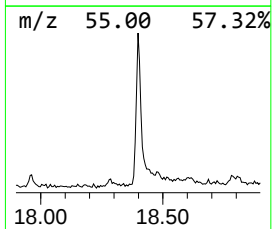
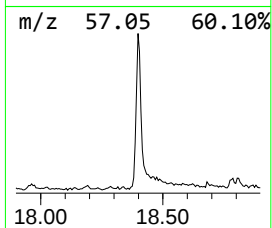
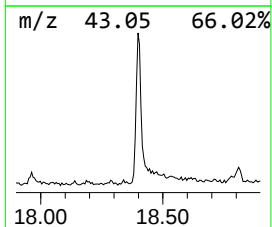
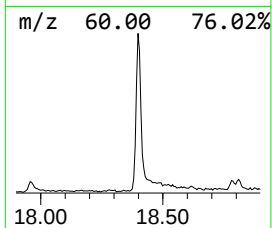
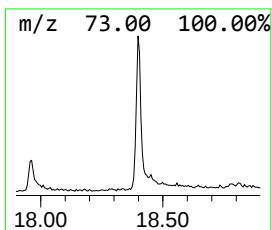
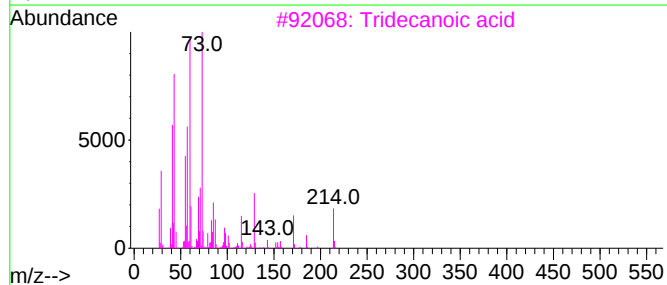
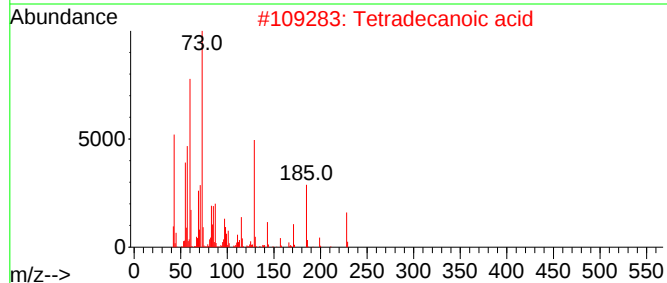
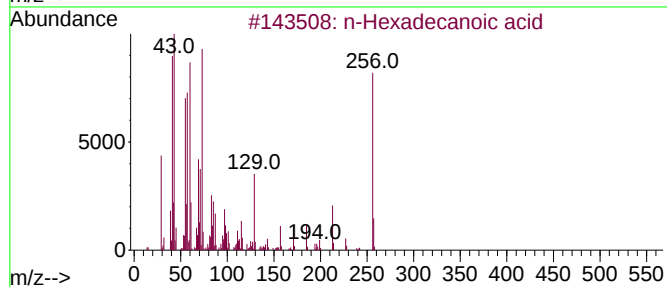
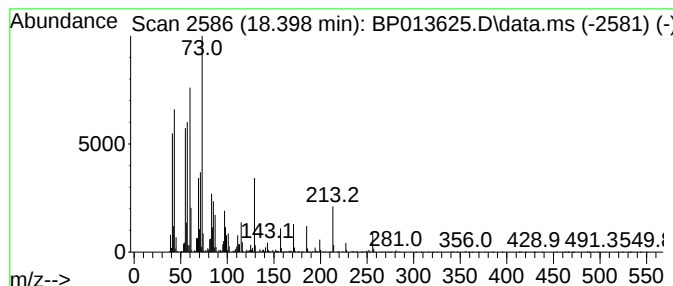
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 n-Hexadecanoic acid Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.398	5.68 ng	677632	Phenanthrene-d10	17.545

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	90
3			Tridecanoic acid	214	C13H26O2	000638-53-9	86
4			Pentadecanoic acid	242	C15H30O2	001002-84-2	81
5			Octadecanoic acid	284	C18H36O2	000057-11-4	81



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP012723\
Data File : BP013625.D
Acq On : 27 Jan 2023 19:24
Operator : CG/JU
Sample : 01320-01
Misc :
ALS Vial : 8 Sample Multiplier: 1

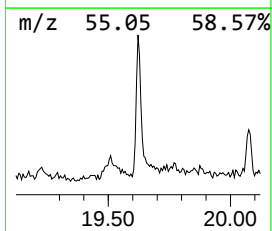
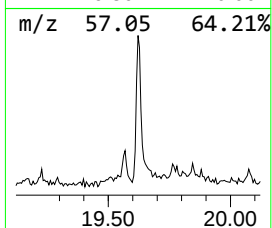
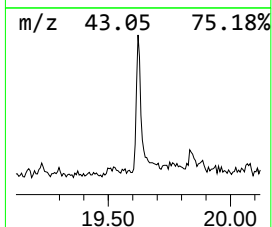
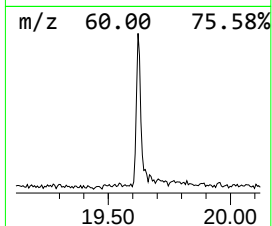
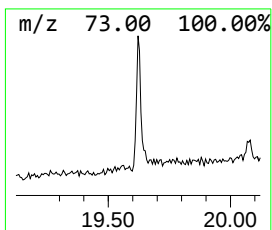
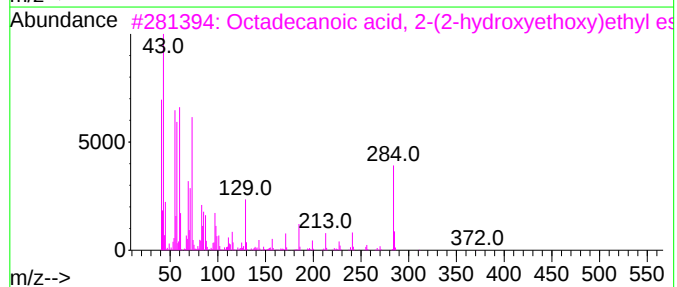
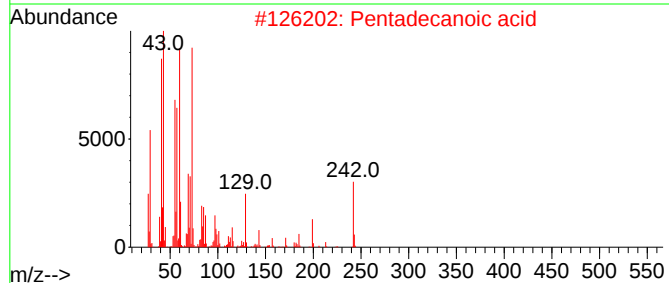
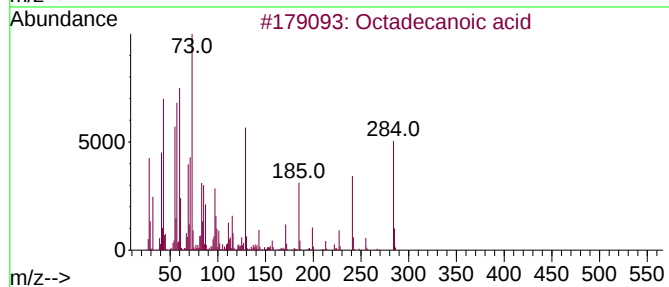
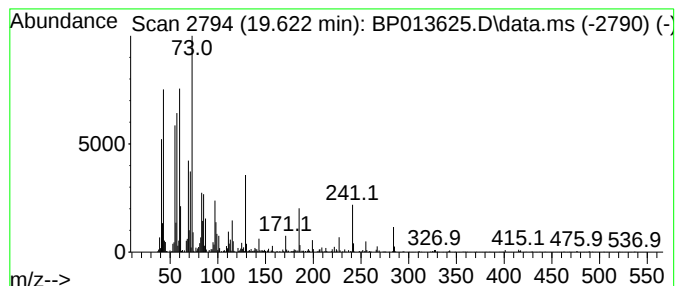
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP012623.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Octadecanoic acid Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.	
19.622	2.05 ng	295968	Chrysene-d12	21.622	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecanoic acid	284	C18H36O2	000057-11-4	99
2	Pentadecanoic acid	242	C15H30O2	001002-84-2	90
3	Octadecanoic acid, 2-(2-hydroxye...	372	C22H44O4	000106-11-6	76
4	Tetradecanoic acid	228	C14H28O2	000544-63-8	74
5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	62



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP012723\
Data File : BP013625.D
Acq On : 27 Jan 2023 19:24
Operator : CG/JU
Sample : 01320-01
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
TOTE-COMP-0126

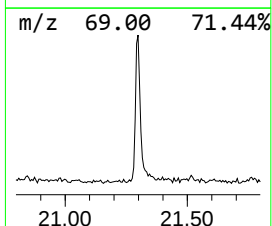
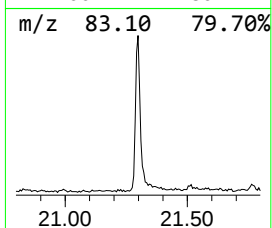
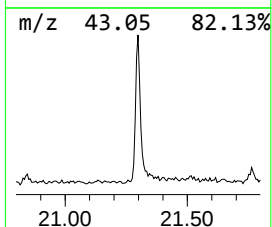
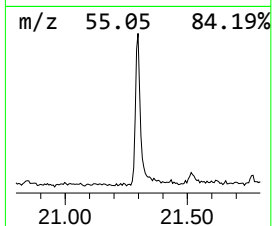
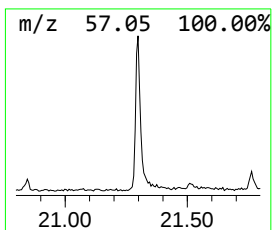
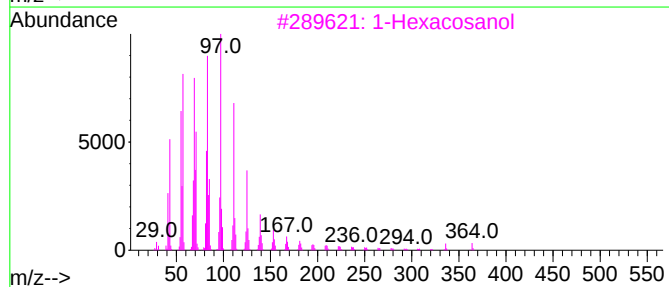
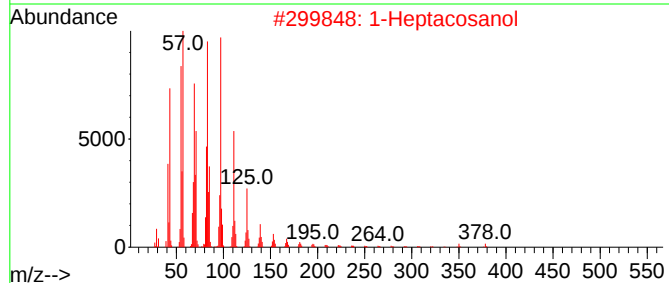
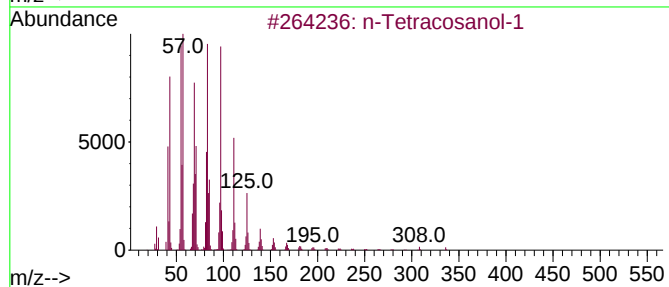
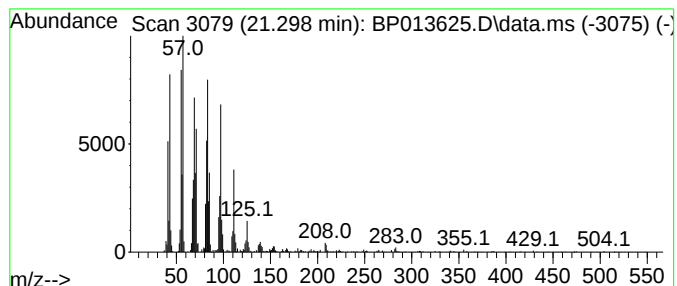
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 n-Tetracosanol-1 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.298	4.25 ng	613063	Chrysene-d12	21.622

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Tetracosanol-1	354	C24H50O	000506-51-4	94
2			1-Heptacosanol	396	C27H56O	002004-39-9	94
3			1-Hexacosanol	382	C26H54O	000506-52-5	91
4			1-Eicosanol	298	C20H42O	000629-96-9	90
5			17-Pentatriacontene	491	C35H70	006971-40-0	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP012723\
Data File : BP013625.D
Acq On : 27 Jan 2023 19:24
Operator : CG/JU
Sample : 01320-01
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
TOTE-COMP-0126

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP012623.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
Butane, 2-metho...	3.223	72.2	ng	3947800	1	8.152	1093250	20.0
Propylene Glycol	3.770	4.3	ng	232913	1	8.152	1093250	20.0
Crotonic acid	4.717	5.8	ng	314532	1	8.152	1093250	20.0
2-Pentanone, 4-...	5.217	13.9	ng	762035	1	8.152	1093250	20.0
Benzene, 1-chlo...	7.117	6.8	ng	368772	1	8.152	1093250	20.0
Benzene, 1-chlo...	7.228	3.1	ng	167369	1	8.152	1093250	20.0
Benzene, (ethox...	8.811	13.9	ng	760149	1	8.152	1093250	20.0
Benzothiazole	11.634	6.0	ng	447521	2	10.981	1478610	20.0
unknown12.581	12.581	13.0	ng	959272	2	10.981	1478610	20.0
Tetradonium Bro...	14.722	5.3	ng	525369	3	14.787	1983150	20.0
Benzophenone	16.145	5.0	ng	494344	3	14.787	1983150	20.0
Oxazolidin-2-on...	17.010	4.9	ng	581843	4	17.545	2385570	20.0
n-Hexadecanoic ...	18.398	5.7	ng	677632	4	17.545	2385570	20.0
Octadecanoic acid	19.622	2.0	ng	295968	5	21.622	2887570	20.0
n-Tetracosanol-1	21.298	4.3	ng	613063	5	21.622	2887570	20.0