

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF030625\
 Data File : BF141867.D
 Acq On : 06 Mar 2025 14:31
 Operator : RC/JU
 Sample : Q1478-01
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 IDW-AQ-MW-19B-COMP-022825

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_F\Methods\8270-BF022725.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF141867.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.163	4	12	27	rBV	3394185	5631353	35.15%	8.430%
2	2.551	62	78	85	rBV	7585521	16022250	100.00%	23.984%
3	3.010	147	156	165	rBV	708771	1346204	8.40%	2.015%
4	3.528	214	244	263	rBV	715035	4601364	28.72%	6.888%
5	4.628	425	431	441	rVB	1753667	2568208	16.03%	3.844%
6	5.098	503	511	521	rBV	133707	205794	1.28%	0.308%
7	5.498	563	579	586	rBV	1931442	2558280	15.97%	3.830%
8	6.439	730	739	744	rBV	124528	195283	1.22%	0.292%
9	6.498	744	749	760	rVB	1248927	1585956	9.90%	2.374%
10	6.886	808	815	821	rBV	680126	925648	5.78%	1.386%
11	7.445	899	910	915	rBV	2596525	3556444	22.20%	5.324%
12	8.063	1010	1015	1022	rBV	258497	336986	2.10%	0.504%
13	8.163	1027	1032	1048	rVB	915768	1177241	7.35%	1.762%
14	9.239	1209	1215	1220	rBV	5053302	6494202	40.53%	9.721%
15	9.916	1325	1330	1335	rBV	1050784	1288901	8.04%	1.929%
16	10.627	1447	1451	1458	rVB	324627	407649	2.54%	0.610%
17	10.704	1458	1464	1469	rBV	3462683	4448109	27.76%	6.658%
18	11.116	1529	1534	1539	rVB	742511	906818	5.66%	1.357%
19	11.398	1577	1582	1595	rVB	1000515	1345288	8.40%	2.014%
20	11.927	1665	1672	1686	rBV2	648109	1036141	6.47%	1.551%
21	12.633	1786	1792	1800	rBV	175522	311847	1.95%	0.467%
22	12.710	1800	1805	1819	rVB	222616	373475	2.33%	0.559%
23	12.986	1846	1852	1858	rBV	5215817	6905363	43.10%	10.337%
24	13.880	2000	2004	2015	rBV2	99740	192560	1.20%	0.288%
25	14.033	2025	2030	2039	rVB	811692	1116705	6.97%	1.672%
26	14.892	2172	2176	2187	rVB	130991	194419	1.21%	0.291%
27	15.509	2275	2281	2295	rVB	652356	1071268	6.69%	1.604%

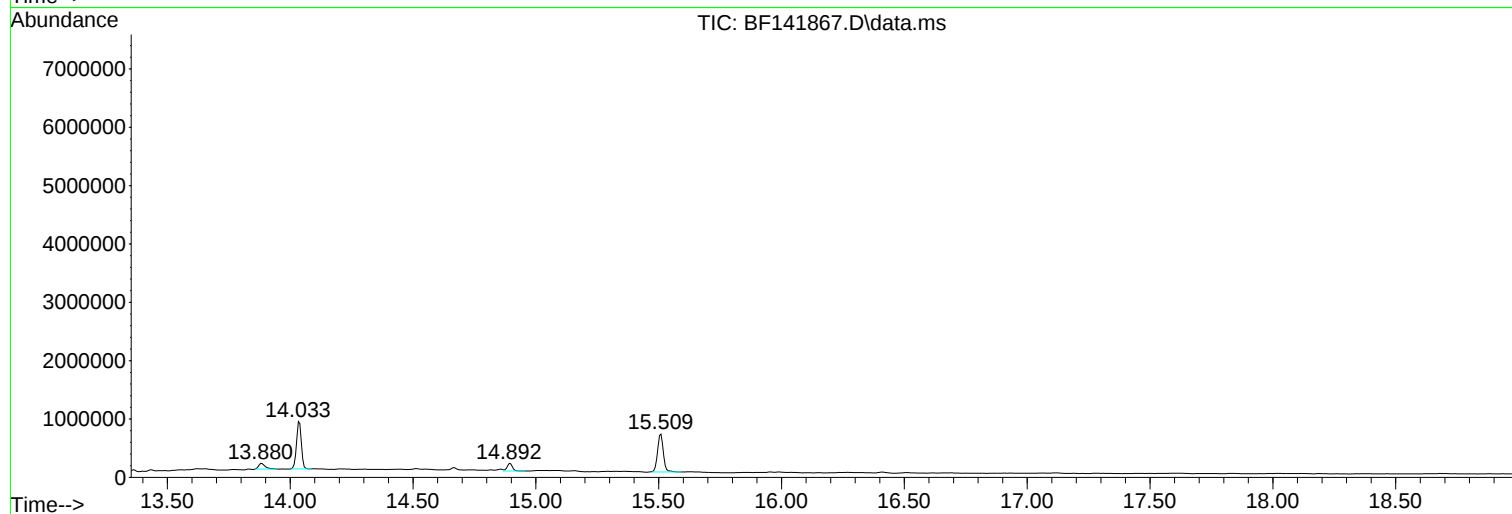
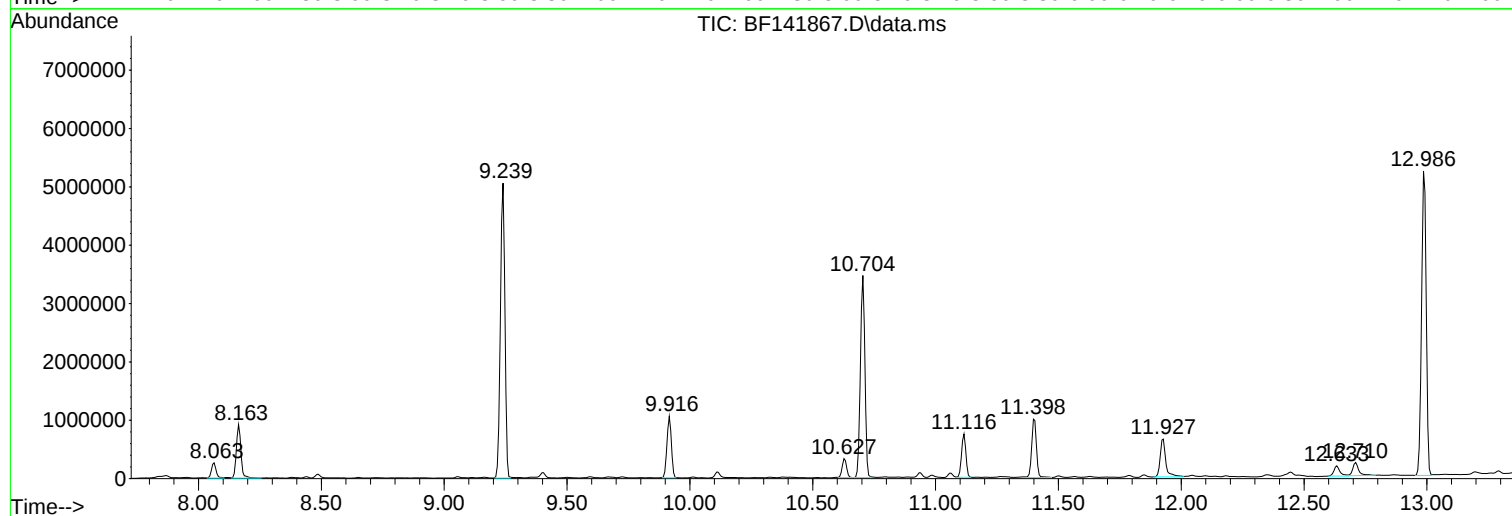
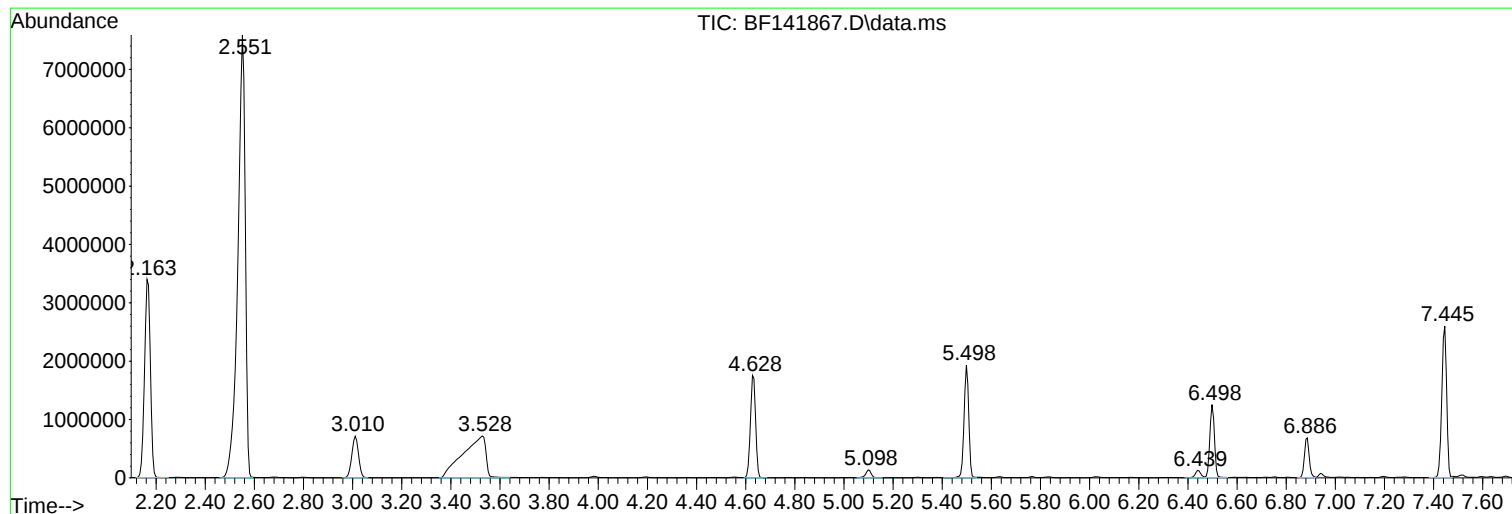
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF022725.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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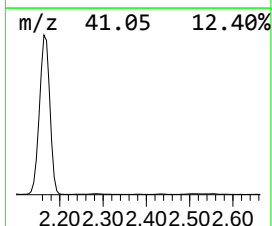
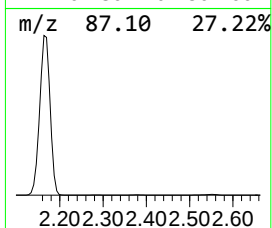
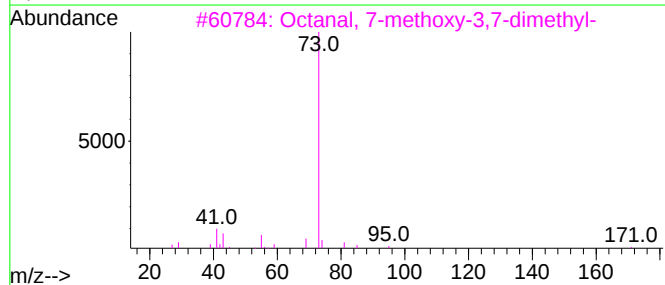
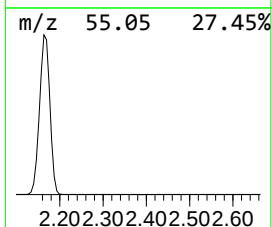
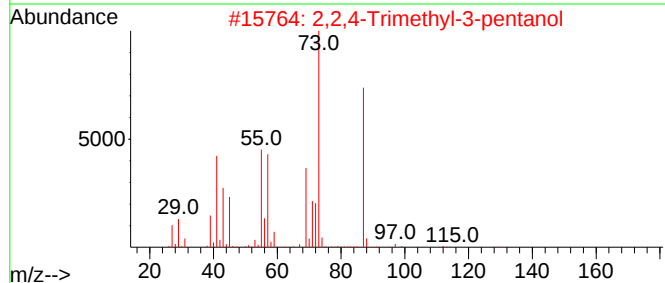
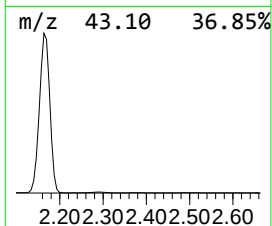
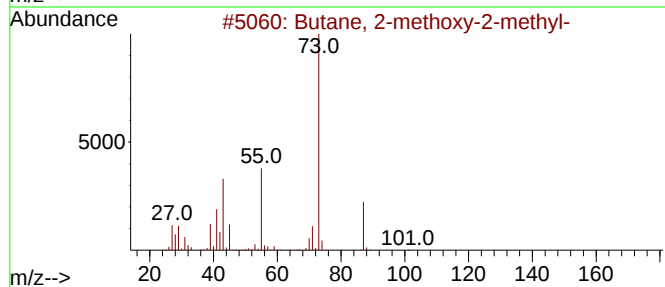
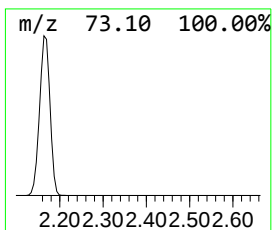
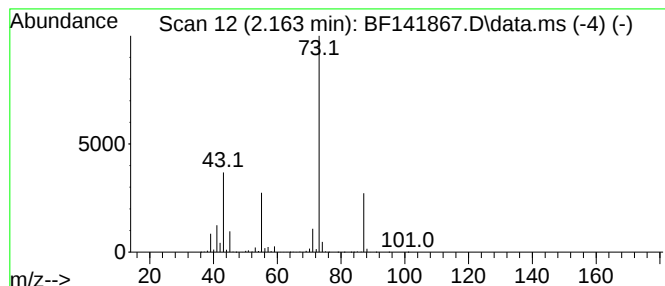
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.163	121.67 ng	5631350	1,4-Dichlorobenzene-d4	6.886

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	40
3			Octanal, 7-methoxy-3,7-dimethyl-	186	C11H22O2	003613-30-7	23
4			Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	22
5			Pentane, 3-methoxy-	102	C6H14O	036839-67-5	17



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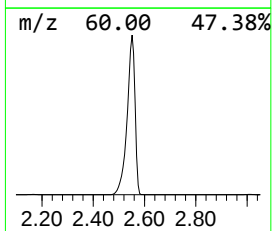
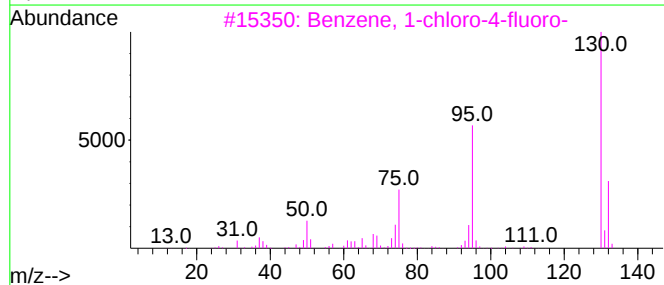
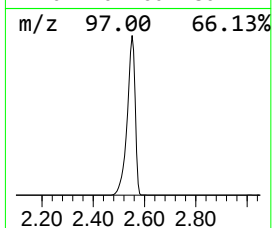
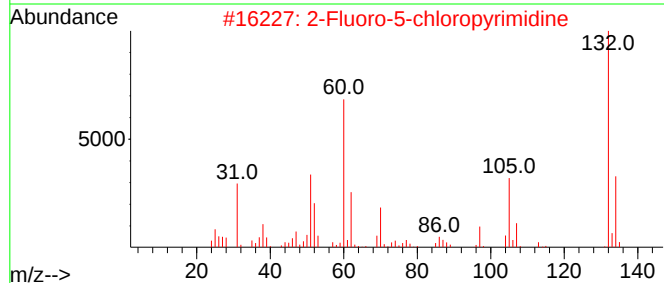
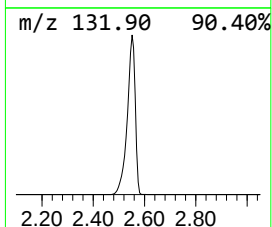
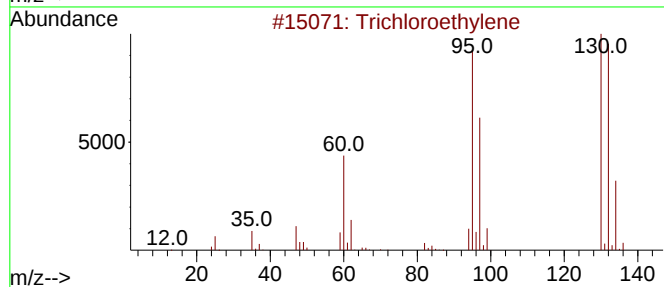
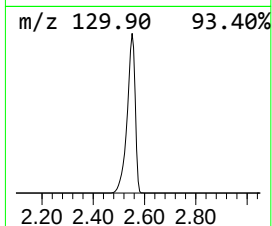
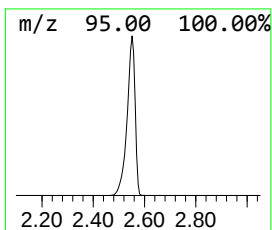
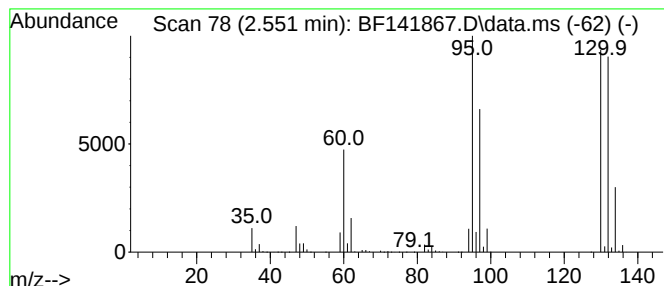
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Trichloroethylene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.551	346.19 ng	16022300	1,4-Dichlorobenzene-d4	6.886

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Trichloroethylene	130	C2HCl3	000079-01-6	99
2			2-Fluoro-5-chloropyrimidine	132	C4H2ClFN2	062802-37-3	35
3			Benzene, 1-chloro-4-fluoro-	130	C6H4ClF	000352-33-0	22
4			Isoquinoline, 3,4-dihydro-	131	C9H9N	003230-65-7	12
5			1H-1,2,4-Triazol-3-amine, 5-(met...	130	C3H6N4S	045534-08-5	10



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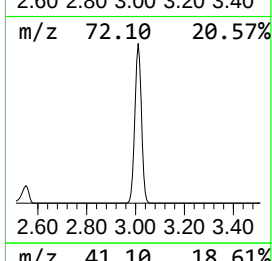
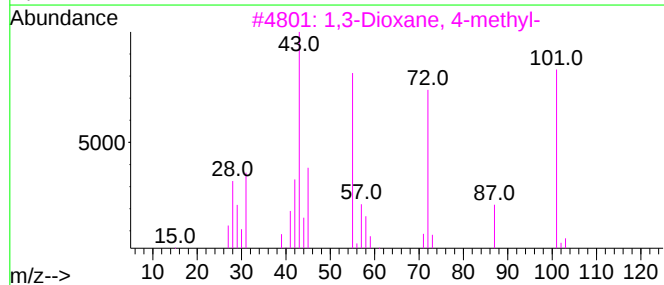
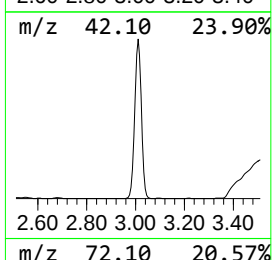
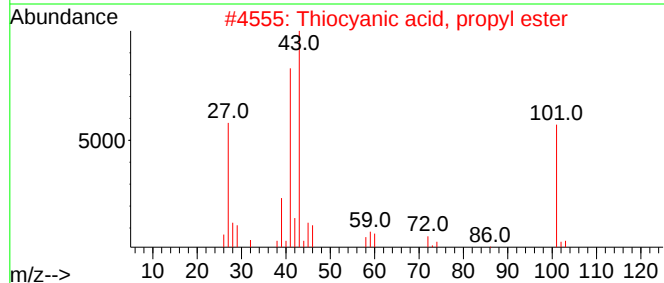
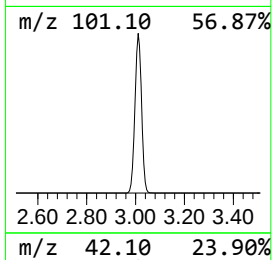
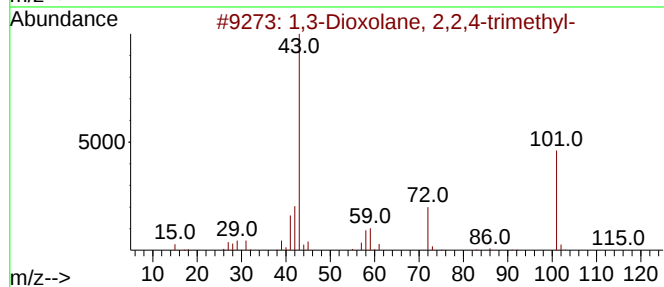
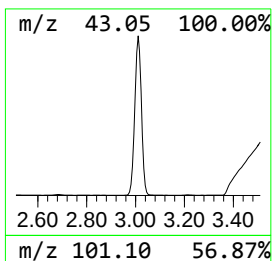
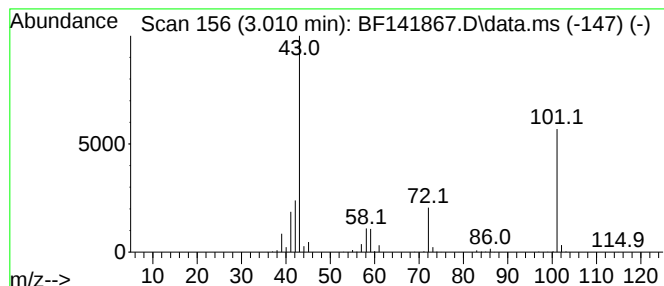
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 1,3-Dioxolane, 2,2,4-trimet... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.010	29.09 ng	1346200	1,4-Dichlorobenzene-d4	6.886

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,3-Dioxolane, 2,2,4-trimethyl-	116	C6H12O2	001193-11-9	90		
2	Thiocyanic acid, propyl ester	101	C4H7NS	004251-16-5	53		
3	1,3-Dioxane, 4-methyl-	102	C5H10O2	001120-97-4	50		
4	Isopropyl isothiocyanate	101	C4H7NS	002253-73-8	50		
5	Propane, 1-isothiocyanato-	101	C4H7NS	000628-30-8	28		



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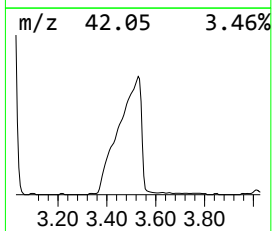
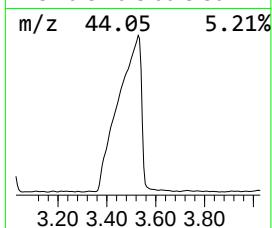
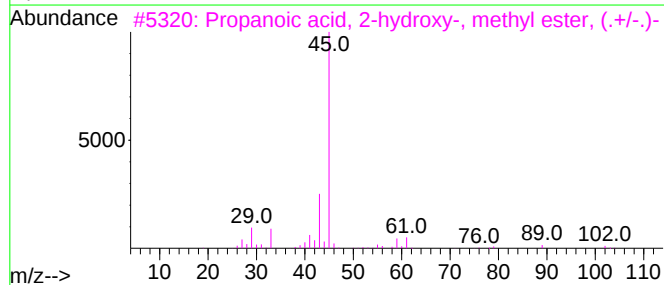
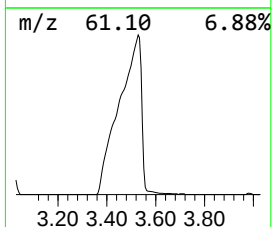
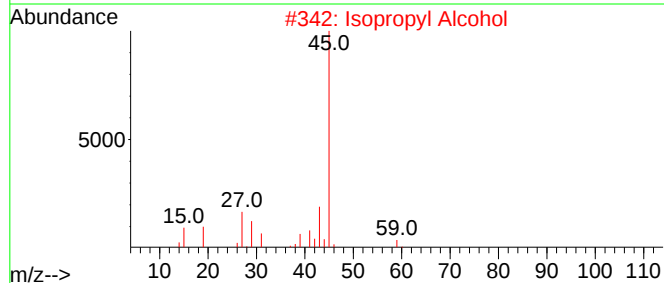
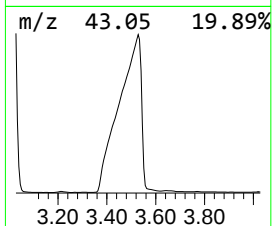
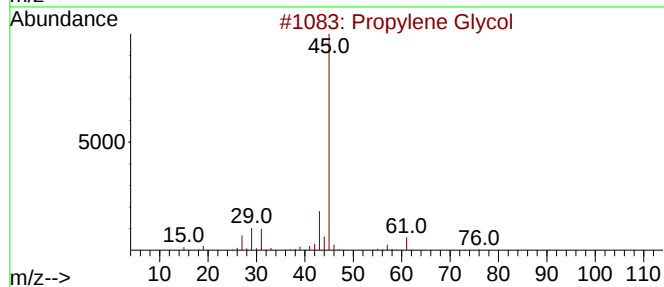
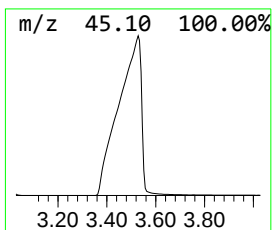
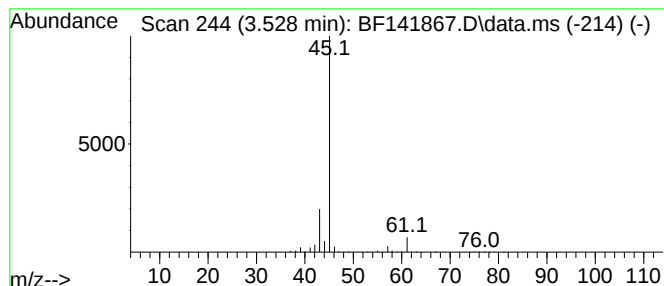
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Peak Number 4 Propylene Glycol Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.528	99.42 ng	4601360	1,4-Dichlorobenzene-d4	6.886

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Propylene Glycol	76	C3H8O2	000057-55-6	90
2			Isopropyl Alcohol	60	C3H8O	000067-63-0	56
3			Propanoic acid, 2-hydroxy-, meth...	104	C4H8O3	000547-64-8	40
4			R-(-)-1,2-propanediol	76	C3H8O2	004254-14-2	9
5			Propane, 2-ethoxy-	88	C5H12O	000625-54-7	9



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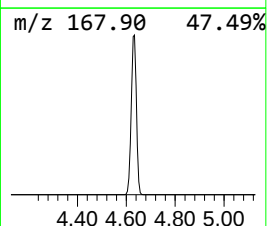
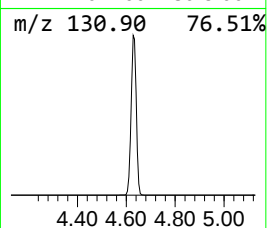
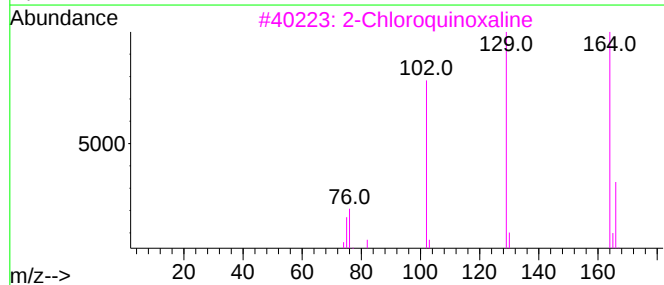
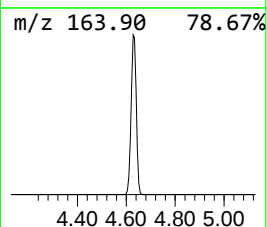
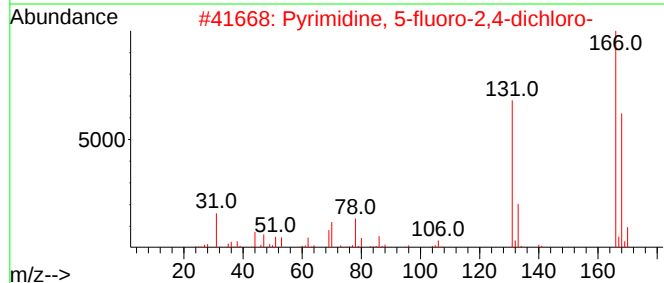
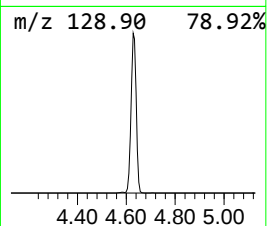
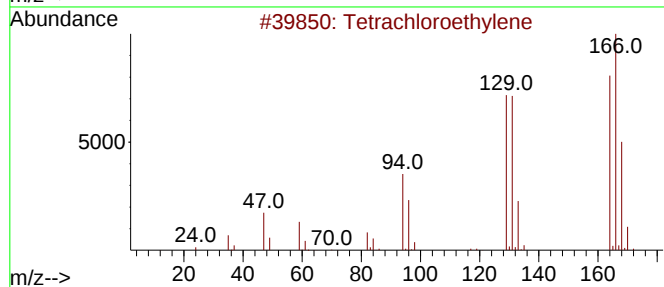
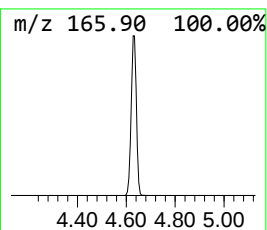
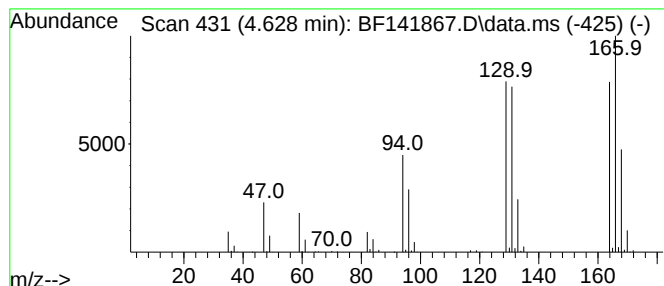
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Tetrachloroethylene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.628	55.49 ng	2568210	1,4-Dichlorobenzene-d4	6.886

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetrachloroethylene	164	C2Cl4	000127-18-4	99
2			Pyrimidine, 5-fluoro-2,4-dichloro-	166	C4HC12FN2	002927-71-1	32
3			2-Chloroquinoxaline	164	C8H5ClN2	001448-87-9	12
4			2,4-Dichloro-6-fluoropyrimidine	166	C4HC12FN2	003833-57-6	9
5			2-Chloro-4-aminopyrimidine	129	C4H4ClN3	007461-50-9	9



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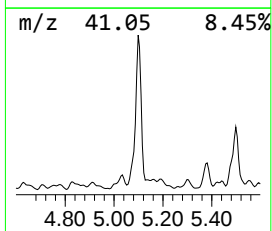
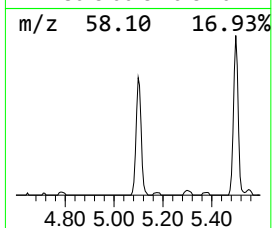
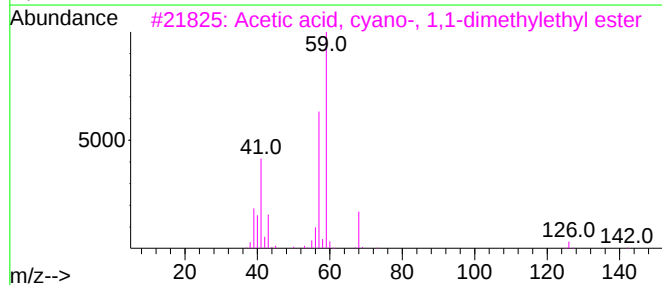
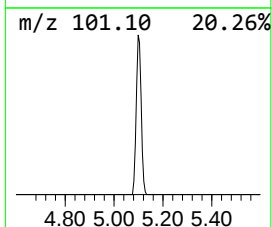
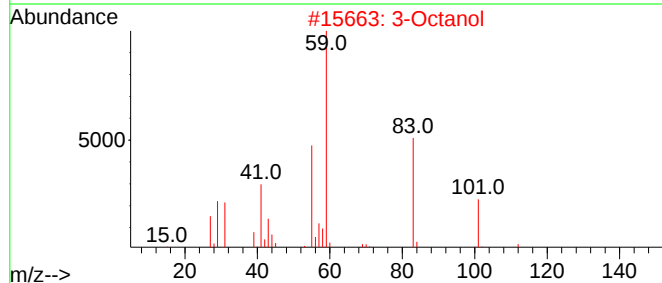
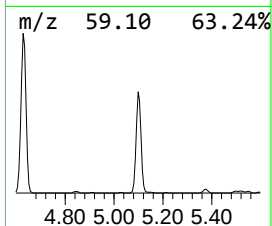
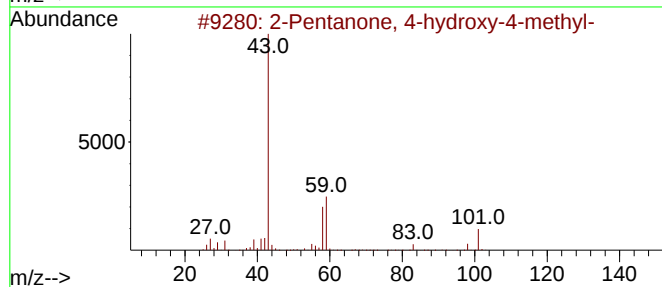
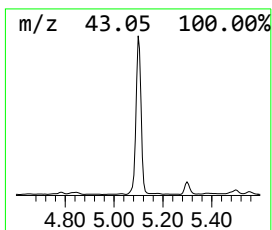
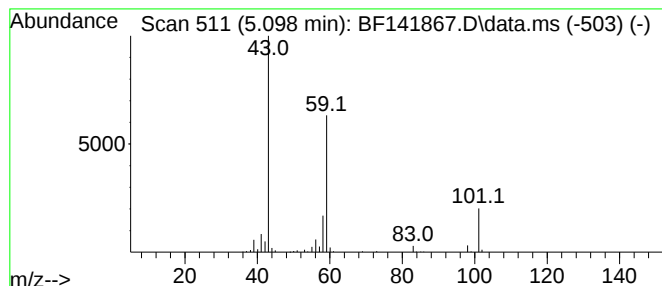
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.098	4.45 ng	205794	1,4-Dichlorobenzene-d4	6.886

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	72
2		3-Octanol	130	C8H18O	000589-98-0	36
3		Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	32
4		1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10
5		Guanidine	59	CH5N3	000113-00-8	9



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF030625\
Data File : BF141867.D
Acq On : 06 Mar 2025 14:31
Operator : RC/JU
Sample : Q1478-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
IDW-AQ-MW-19B-COMP-022825

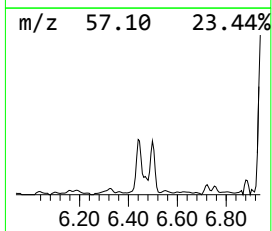
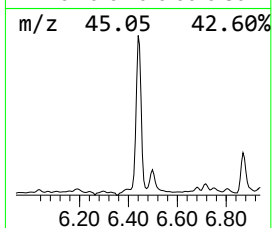
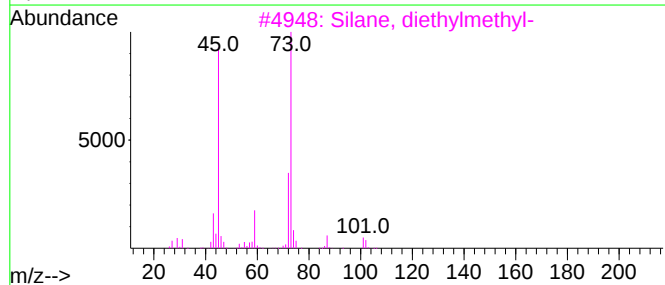
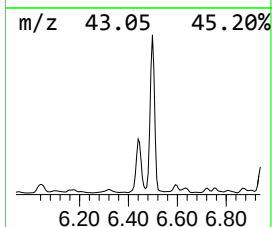
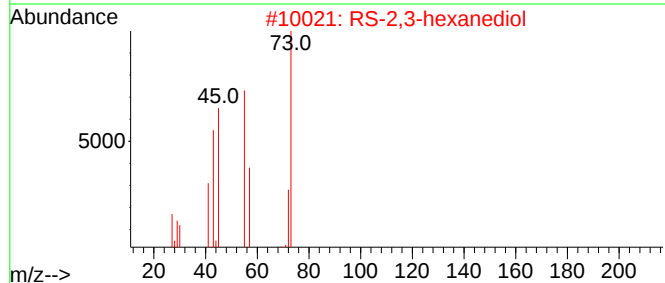
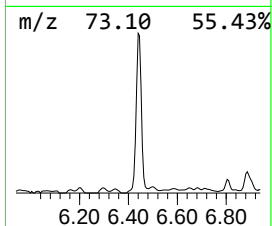
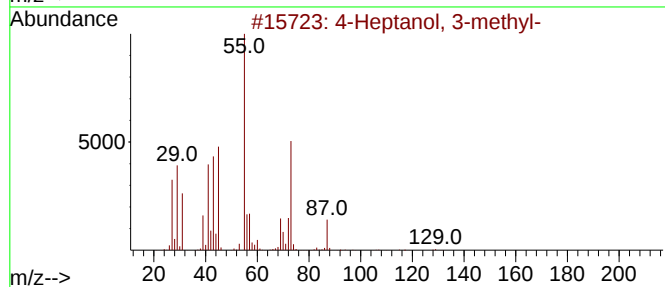
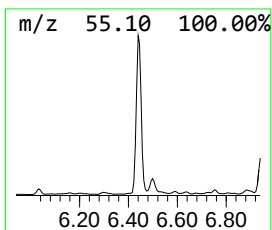
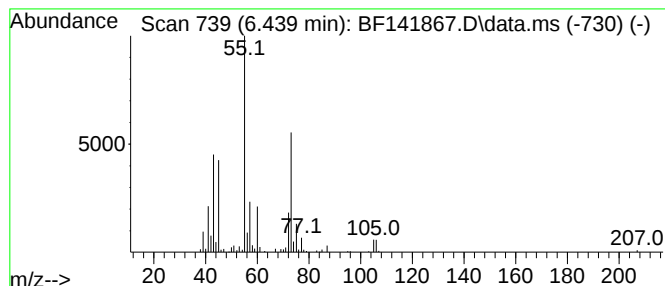
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 unknown6.439 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.439	4.22 ng	195283	1,4-Dichlorobenzene-d4	6.886

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4-Heptanol, 3-methyl-	130	C8H18O	001838-73-9	40
2			RS-2,3-hexanediol	118	C6H14O2	082360-67-6	38
3			Silane, diethylmethyl-	102	C5H14Si	000760-32-7	32
4			3-Hexanol, 2-methyl-	116	C7H16O	000617-29-8	27
5			5-Hydroxy-4-octanone	144	C8H16O2	000496-77-5	27



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF030625\
Data File : BF141867.D
Acq On : 06 Mar 2025 14:31
Operator : RC/JU
Sample : Q1478-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
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ClientSampleId :
IDW-AQ-MW-19B-COMP-022825

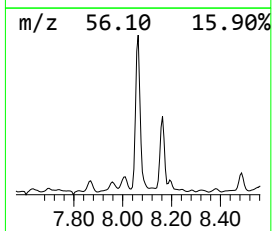
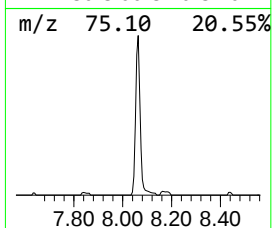
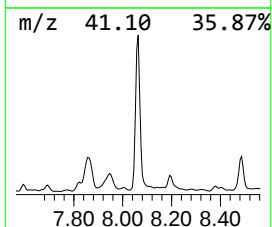
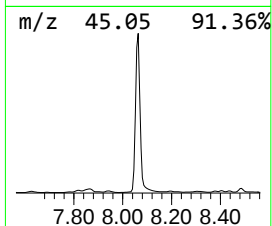
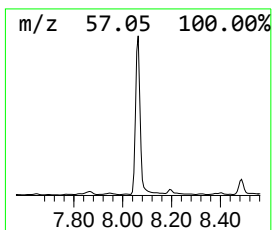
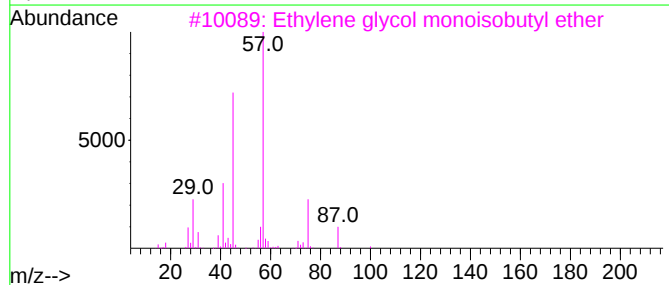
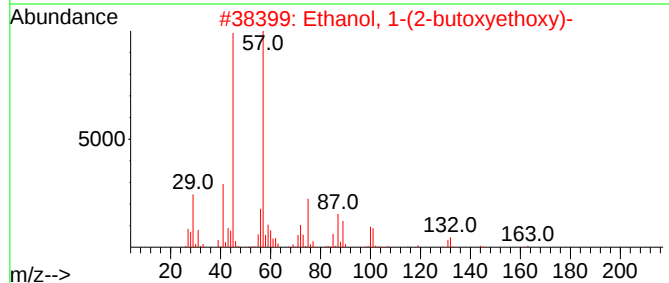
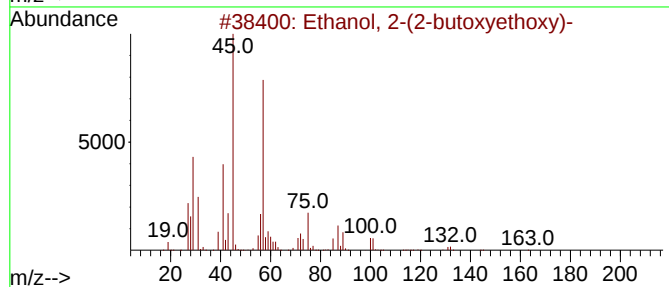
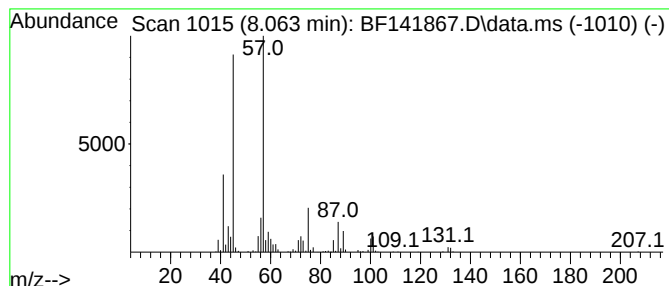
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Ethanol, 2-(2-butoxyethoxy)- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.063	5.73 ng	336986	Naphthalene-d8	8.163

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	83
3			Ethylene glycol monoisobutyl ether	118	C6H14O2	004439-24-1	64
4			Ethanol, 2-[2-(2-methylpropoxy)e...	162	C8H18O3	018912-80-6	56
5			3,3-Dimethylbutane-2-ol	102	C6H14O	000464-07-3	52



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF030625\
Data File : BF141867.D
Acq On : 06 Mar 2025 14:31
Operator : RC/JU
Sample : Q1478-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

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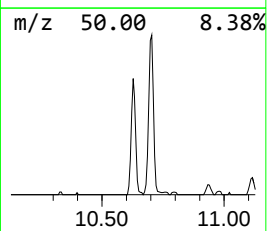
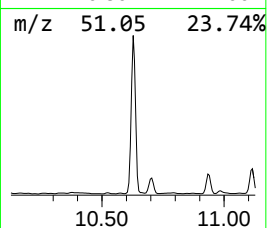
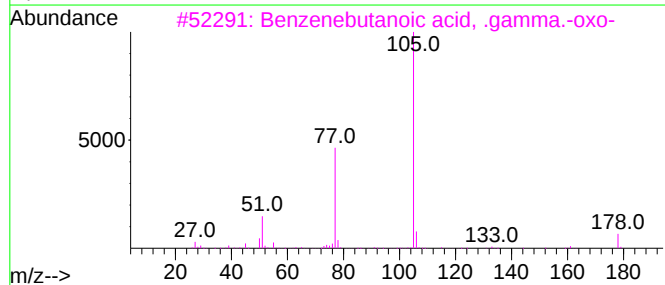
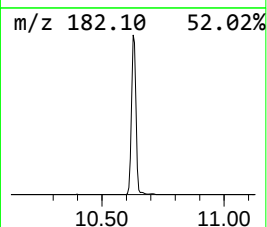
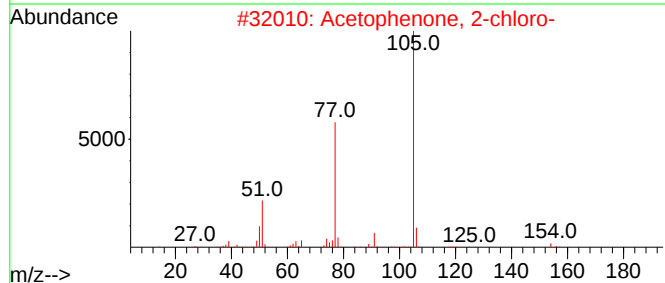
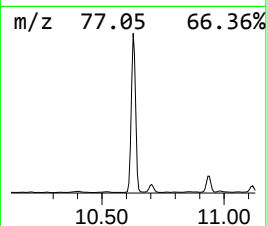
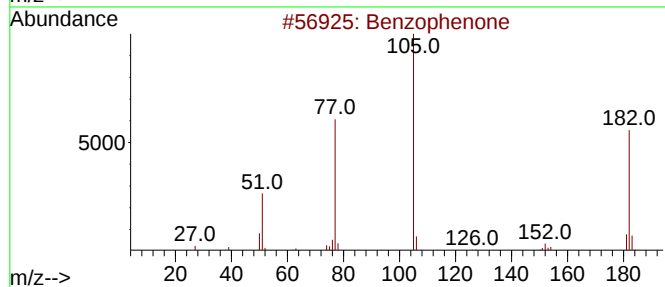
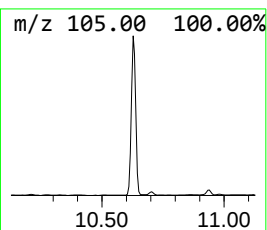
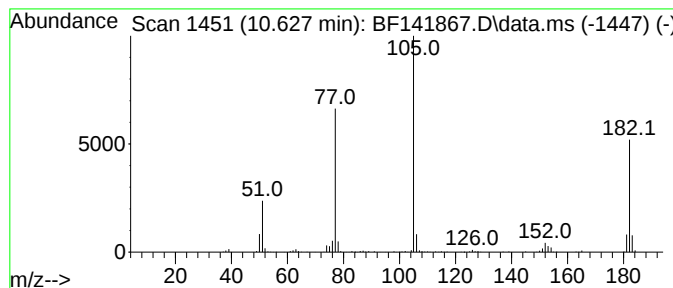
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Benzophenone Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.627	6.33 ng	407649	Acenaphthene-d10	9.916

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	96
2			Acetophenone, 2-chloro-	154	C8H7ClO	000532-27-4	49
3			Benzenebutanoic acid, .gamma.-oxo-	178	C10H10O3	002051-95-8	47
4			N-(1-Cyanovinyl)benzamide	172	C10H8N2O	1000186-19-1	47
5			Vinyl benzoate	148	C9H8O2	000769-78-8	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF030625\
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Operator : RC/JU
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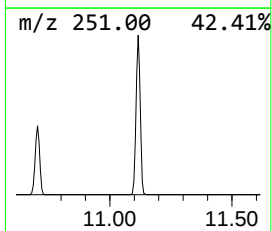
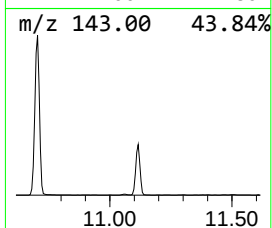
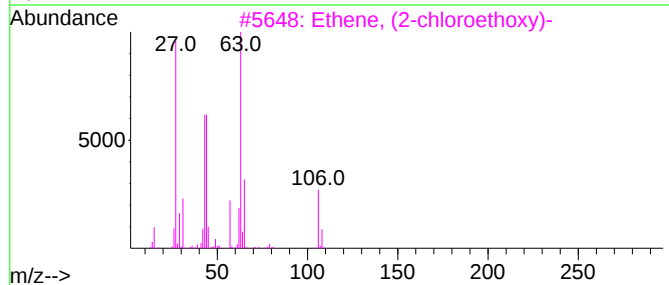
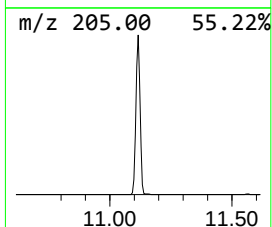
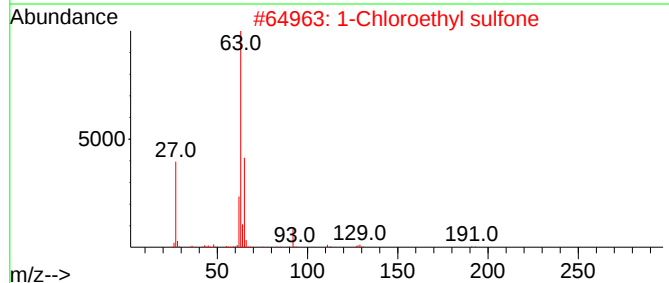
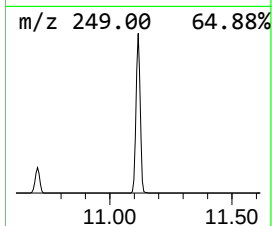
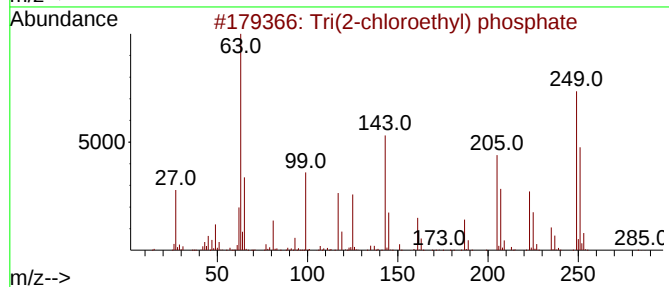
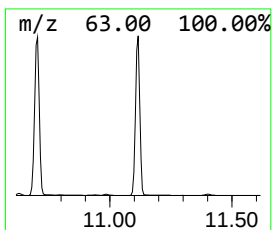
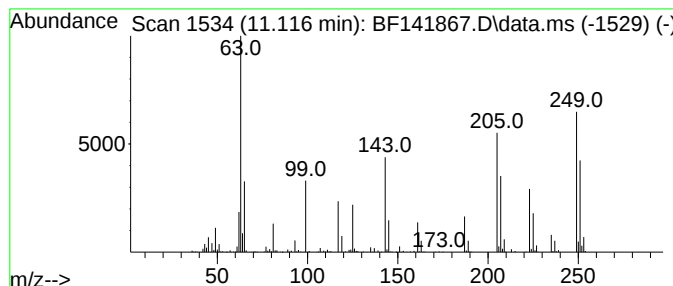
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST0.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 Tri(2-chloroethyl) phosphate Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.116	13.48 ng	906818	Phenanthrene-d10	11.398

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tri(2-chloroethyl) phosphate	284	C6H12Cl3O4P	000115-96-8	95
2			1-Chloroethyl sulfone	190	C4H8Cl2O2S	010038-07-0	47
3			Ethene, (2-chloroethoxy)-	106	C4H7ClO	000110-75-8	47
4			Ethanol, 2-chloro-, phosphite (3:1)	268	C6H12Cl3O3P	000140-08-9	38
5			Trichloroacetic acid, 2-chloroet...	224	C4H4Cl4O2	1000330-88-7	37



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF030625\
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Acq On : 06 Mar 2025 14:31
Operator : RC/JU
Sample : Q1478-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
IDW-AQ-MW-19B-COMP-022825

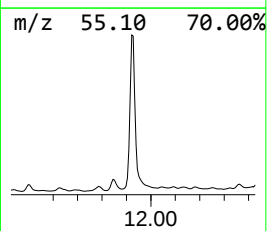
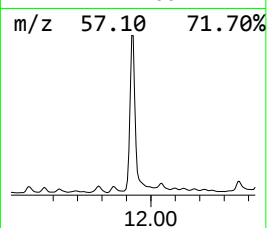
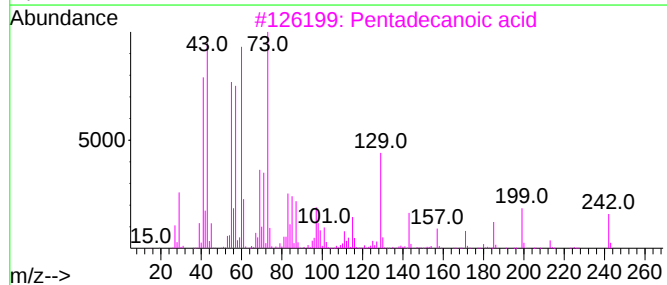
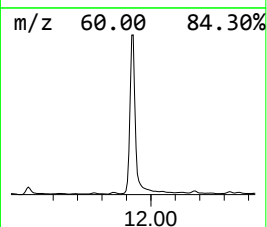
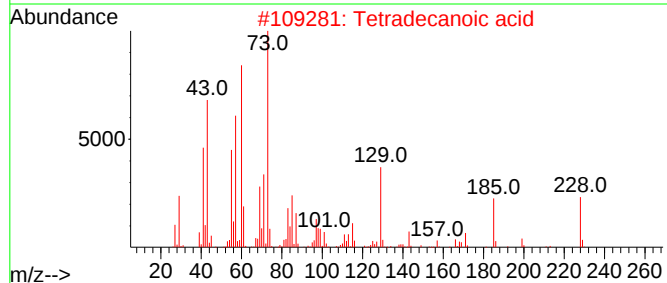
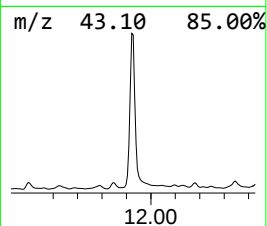
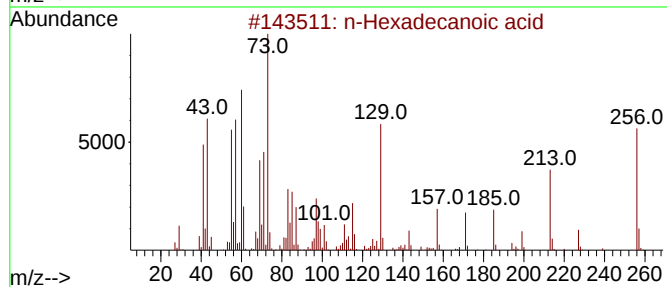
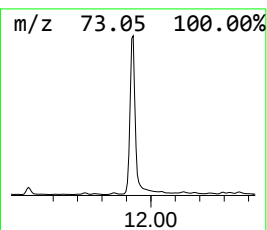
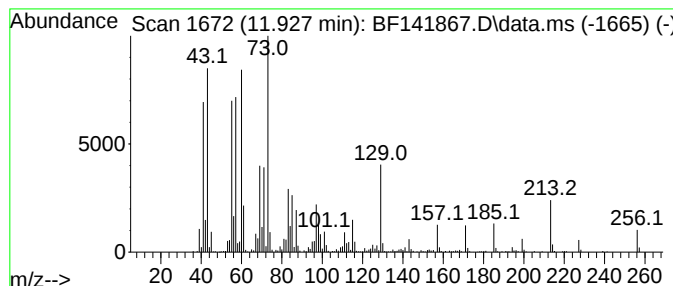
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 n-Hexadecanoic acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.927	15.40 ng	1036140	Phenanthrene-d10	11.398

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	94
3			Pentadecanoic acid	242	C15H30O2	001002-84-2	93
4			Tridecanoic acid	214	C13H26O2	000638-53-9	91
5			n-Decanoic acid	172	C10H20O2	000334-48-5	68



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF030625\
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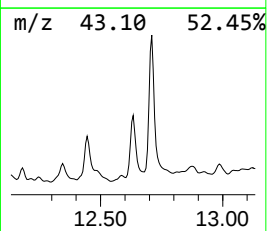
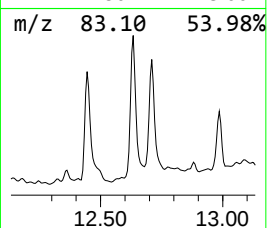
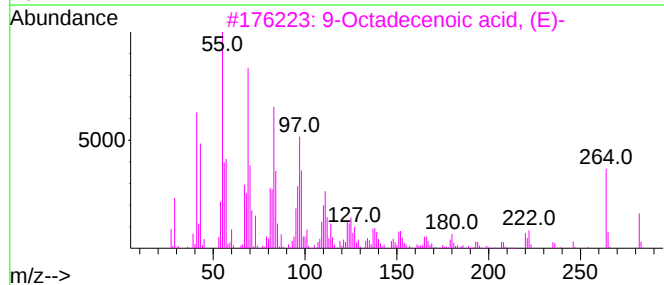
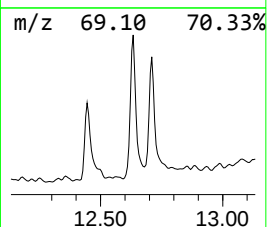
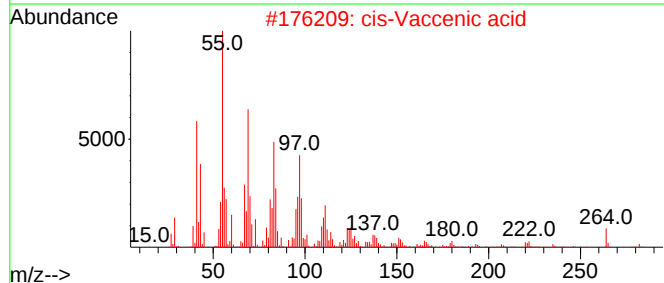
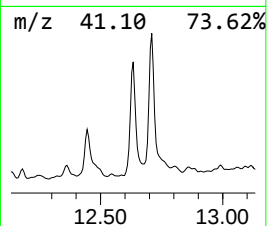
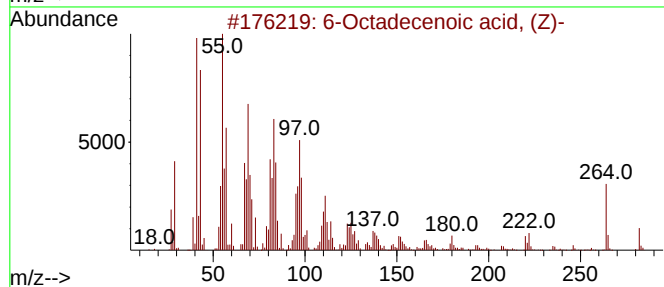
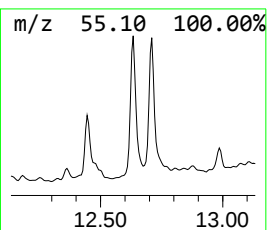
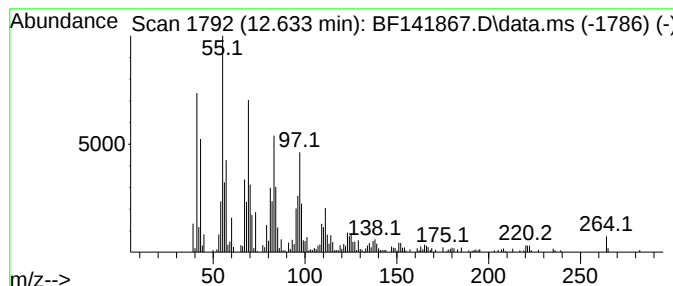
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 6-Octadecenoic acid, (Z)- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.633	4.64 ng	311847	Phenanthrene-d10	11.398

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			6-Octadecenoic acid, (Z)-	282	C18H34O2	000593-39-5	99
2			cis-Vaccenic acid	282	C18H34O2	000506-17-2	99
3			9-Octadecenoic acid, (E)-	282	C18H34O2	000112-79-8	99
4			6-Octadecenoic acid	282	C18H34O2	1000336-66-8	99
5			trans-13-Octadecenoic acid	282	C18H34O2	000693-71-0	98



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF030625\
Data File : BF141867.D
Acq On : 06 Mar 2025 14:31
Operator : RC/JU
Sample : Q1478-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
IDW-AQ-MW-19B-COMP-022825

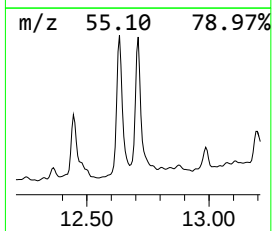
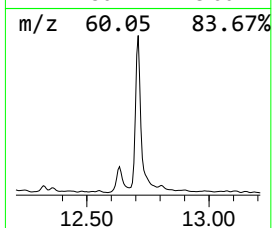
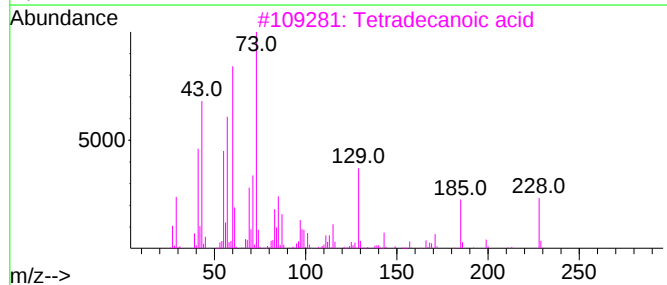
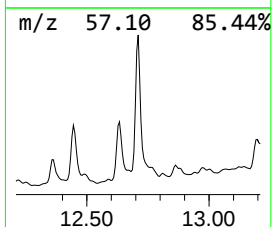
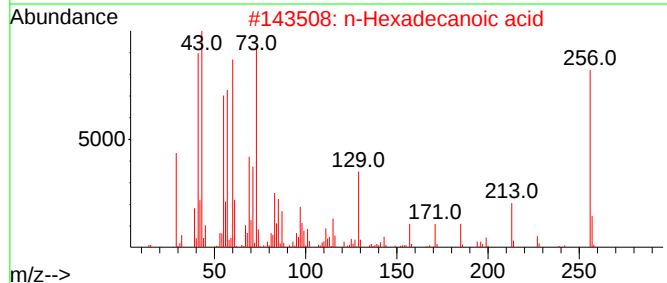
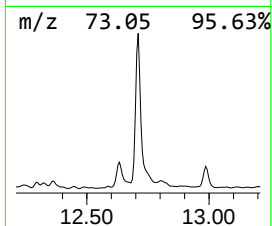
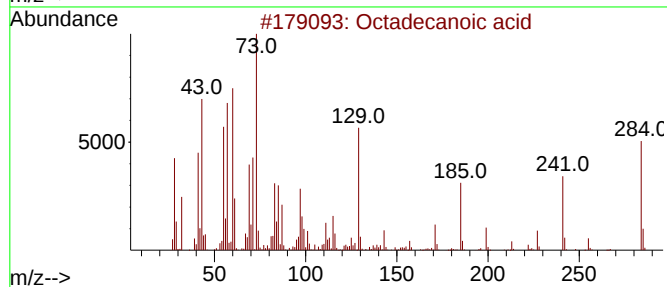
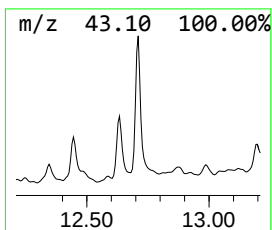
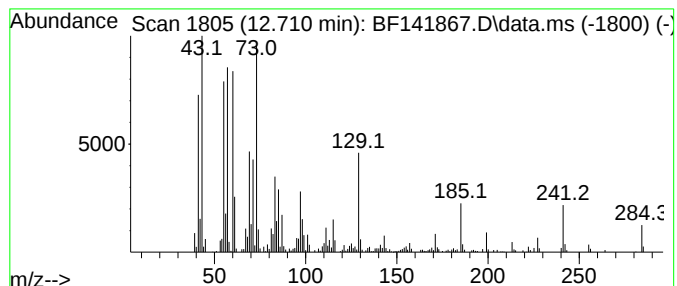
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Octadecanoic acid Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.710	5.55 ng	373475	Phenanthrene-d10	11.398

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid	284	C18H36O2	000057-11-4	99
2			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	91
3			Tetradecanoic acid	228	C14H28O2	000544-63-8	83
4			Pentadecanoic acid	242	C15H30O2	001002-84-2	72
5			Tridecanoic acid	214	C13H26O2	000638-53-9	62



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF030625\
Data File : BF141867.D
Acq On : 06 Mar 2025 14:31
Operator : RC/JU
Sample : Q1478-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

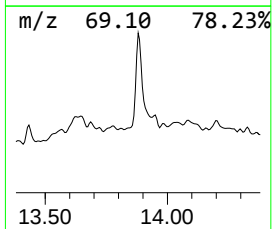
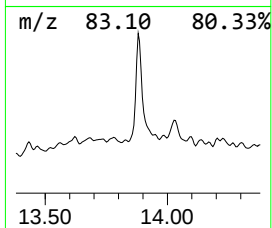
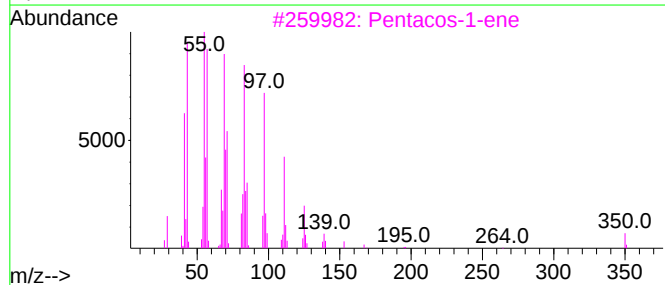
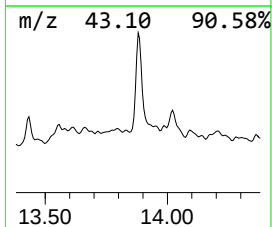
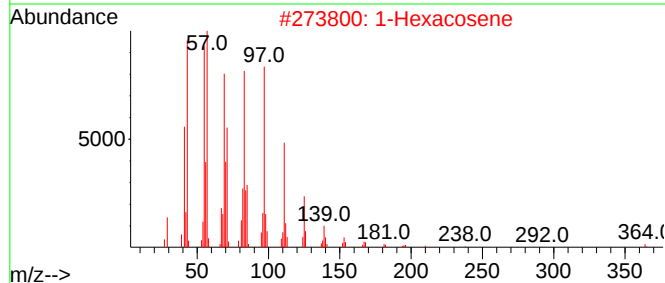
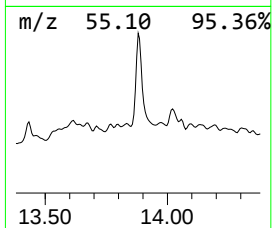
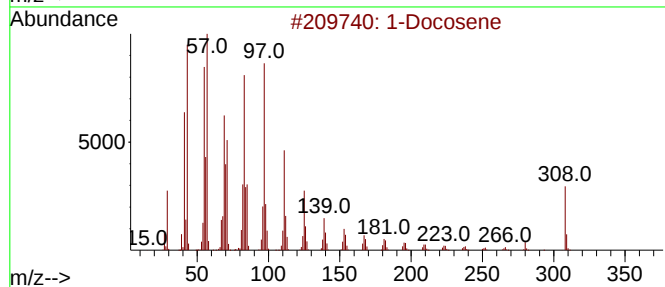
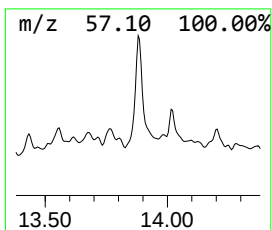
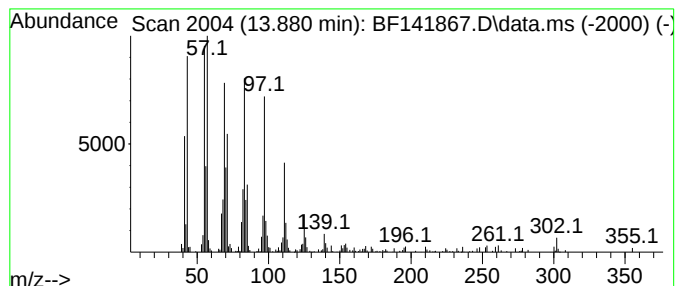
Instrument :
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ClientSampleId :
IDW-AQ-MW-19B-COMP-022825

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF022725.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 1-Docosene Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.880	3.45 ng	192560	Chrysene-d12	14.033	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Docosene	308	C22H44	001599-67-3	95
2	1-Hexacosene	364	C26H52	018835-33-1	95
3	Pentacos-1-ene	350	C25H50	016980-85-1	94
4	Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	94
5	1-Heneicosyl formate	340	C22H44O2	077899-03-7	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF030625\
Data File : BF141867.D
Acq On : 06 Mar 2025 14:31
Operator : RC/JU
Sample : Q1478-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
IDW-AQ-MW-19B-COMP-022825

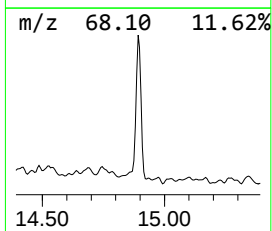
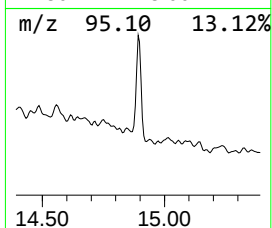
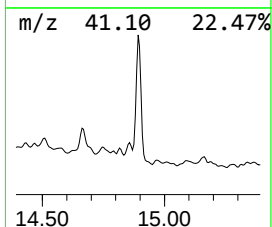
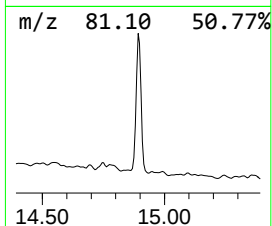
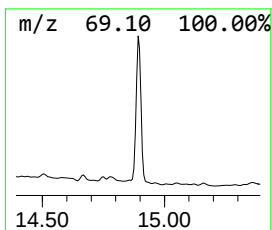
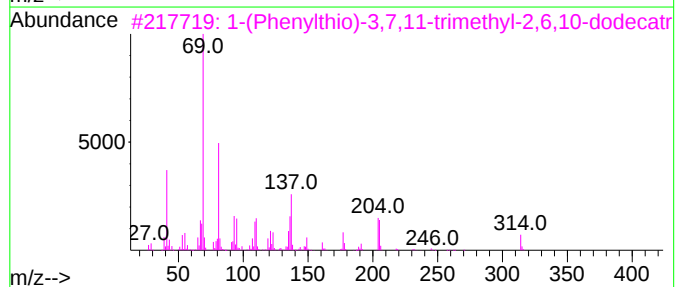
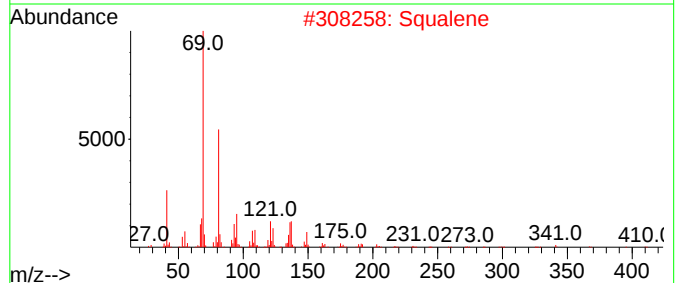
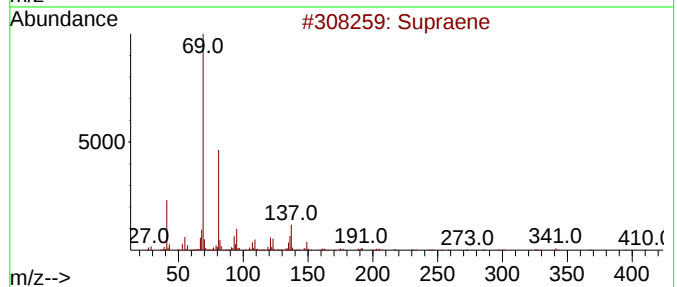
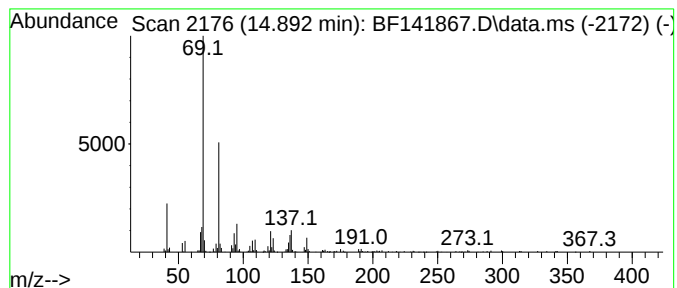
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF022725.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Supraene Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.892	3.63 ng	194419	Perylene-d12	15.509

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Supraene	410	C30H50	007683-64-9	91
2			Squalene	410	C30H50	000111-02-4	90
3			1-(Phenylthio)-3,7,11-trimethyl-...	314	C21H30S	028413-57-2	87
4			4,8,12-Tetradecatrienal, 5,9,13-...	248	C17H28O	066408-55-7	72
5			2,6,10,14-Hexadecatetraen-1-ol, ...	332	C22H36O2	061691-98-3	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF030625\
Data File : BF141867.D
Acq On : 06 Mar 2025 14:31
Operator : RC/JU
Sample : Q1478-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
IDW-AQ-MW-19B-COMP-022825

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF022725.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...	2.163	121.7	ng	5631350	1	6.886	925648	20.0
Trichloroethylene	2.551	346.2	ng	16022300	1	6.886	925648	20.0
1,3-Dioxolane, ...	3.010	29.1	ng	1346200	1	6.886	925648	20.0
Propylene Glycol	3.528	99.4	ng	4601360	1	6.886	925648	20.0
Tetrachloroethy...	4.628	55.5	ng	2568210	1	6.886	925648	20.0
2-Pentanone, 4-...	5.098	4.5	ng	205794	1	6.886	925648	20.0
unknown6.439	6.439	4.2	ng	195283	1	6.886	925648	20.0
Ethanol, 2-(2-b...	8.063	5.7	ng	336986	2	8.163	1177240	20.0
Benzophenone	10.627	6.3	ng	407649	3	9.916	1288900	20.0
Tri(2-chloroeth...	11.116	13.5	ng	906818	4	11.398	1345290	20.0
n-Hexadecanoic ...	11.927	15.4	ng	1036140	4	11.398	1345290	20.0
6-Octadecenoic ...	12.633	4.6	ng	311847	4	11.398	1345290	20.0
Octadecanoic acid	12.710	5.5	ng	373475	4	11.398	1345290	20.0
1-Docosene	13.880	3.5	ng	192560	5	14.033	1116710	20.0
Supraene	14.892	3.6	ng	194419	6	15.509	1071270	20.0