

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF111022\
 Data File : BF131117.D
 Acq On : 10 Nov 2022 11:51
 Operator : CG\JU
 Sample : N5495-10
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 DUP-01

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-BF110922.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF131117.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.199	11	19	27	rVB	327537	442270	23.52%	3.155%
2	2.310	33	38	42	rVB	14377	19364	1.03%	0.138%
3	3.028	154	160	167	rVB	42100	66514	3.54%	0.475%
4	4.498	406	410	415	rVB	30984	33706	1.79%	0.240%
5	5.116	510	515	526	rVB	286546	357856	19.03%	2.553%
6	5.510	578	582	586	rBV	1855644	1880797	100.00%	13.418%
7	6.516	748	753	756	rBV	1796175	1826120	97.09%	13.028%
8	6.887	812	816	819	rBV	492999	354910	18.87%	2.532%
9	7.451	907	912	915	rBV	1059944	1055886	56.14%	7.533%
10	8.169	1030	1034	1038	rBV	461036	444944	23.66%	3.174%
11	9.251	1212	1218	1221	rBV	1972943	1838635	97.76%	13.117%
12	9.933	1329	1334	1337	rBV	577163	503208	26.76%	3.590%
13	10.727	1464	1469	1472	rBV	1256678	1099520	58.46%	7.844%
14	11.427	1584	1588	1591	rBV	640790	515044	27.38%	3.674%
15	12.392	1739	1752	1764	rBV6	26515	86721	4.61%	0.619%
16	12.516	1768	1773	1778	rBV	20877	27034	1.44%	0.193%
17	13.016	1854	1858	1861	rBV	1872724	1605122	85.34%	11.451%
18	13.392	1913	1922	1931	rBV2	44732	101734	5.41%	0.726%
19	13.574	1949	1953	1956	rVB3	14706	20835	1.11%	0.149%
20	13.980	2017	2022	2030	rBV3	44160	96834	5.15%	0.691%
21	14.074	2034	2038	2048	rVB	416931	423381	22.51%	3.020%
22	14.168	2050	2054	2059	rVV	50817	52765	2.81%	0.376%
23	14.657	2132	2137	2142	rVV	49661	75809	4.03%	0.541%
24	14.710	2143	2146	2153	rVB8	16313	25343	1.35%	0.181%
25	14.845	2163	2169	2176	rVB5	26556	53817	2.86%	0.384%
26	14.921	2179	2182	2186	rVB	22246	24337	1.29%	0.174%
27	15.239	2231	2236	2241	rVB3	31825	56921	3.03%	0.406%
28	15.574	2288	2293	2299	rVB	260031	328318	17.46%	2.342%
29	15.898	2344	2348	2358	rVB4	27133	58640	3.12%	0.418%
30	16.421	2426	2437	2458	rBV4	69134	384493	20.44%	2.743%
31	17.086	2544	2550	2558	rVB10	21796	53278	2.83%	0.380%
32	18.668	2811	2819	2829	rVB6	30733	103157	5.48%	0.736%

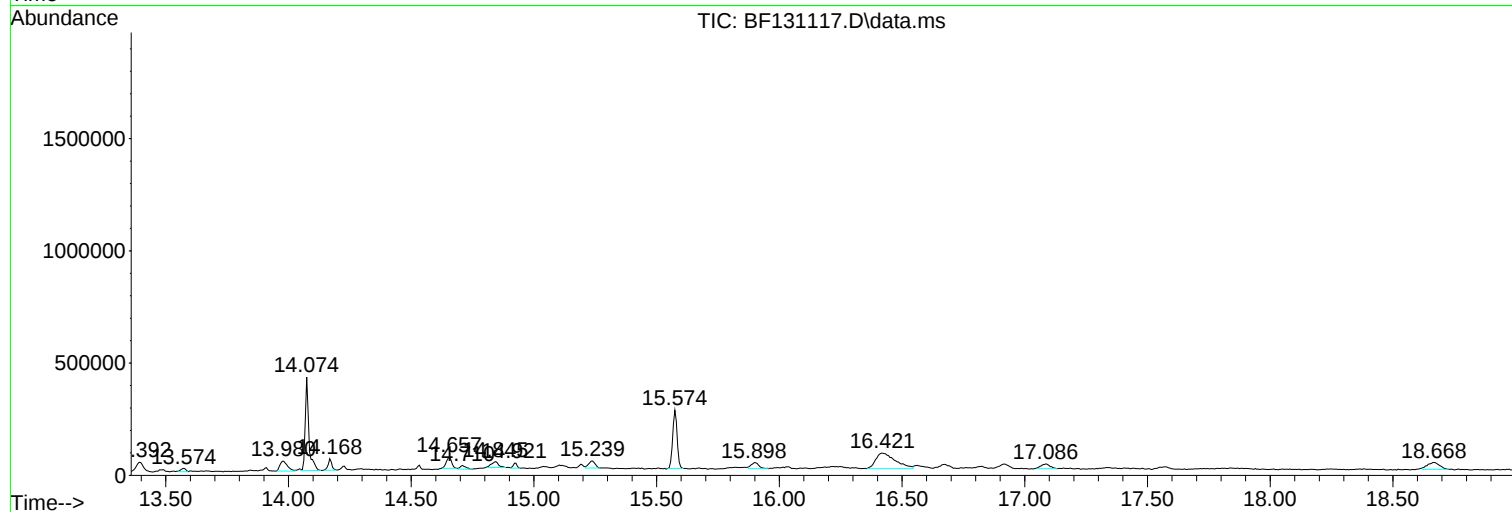
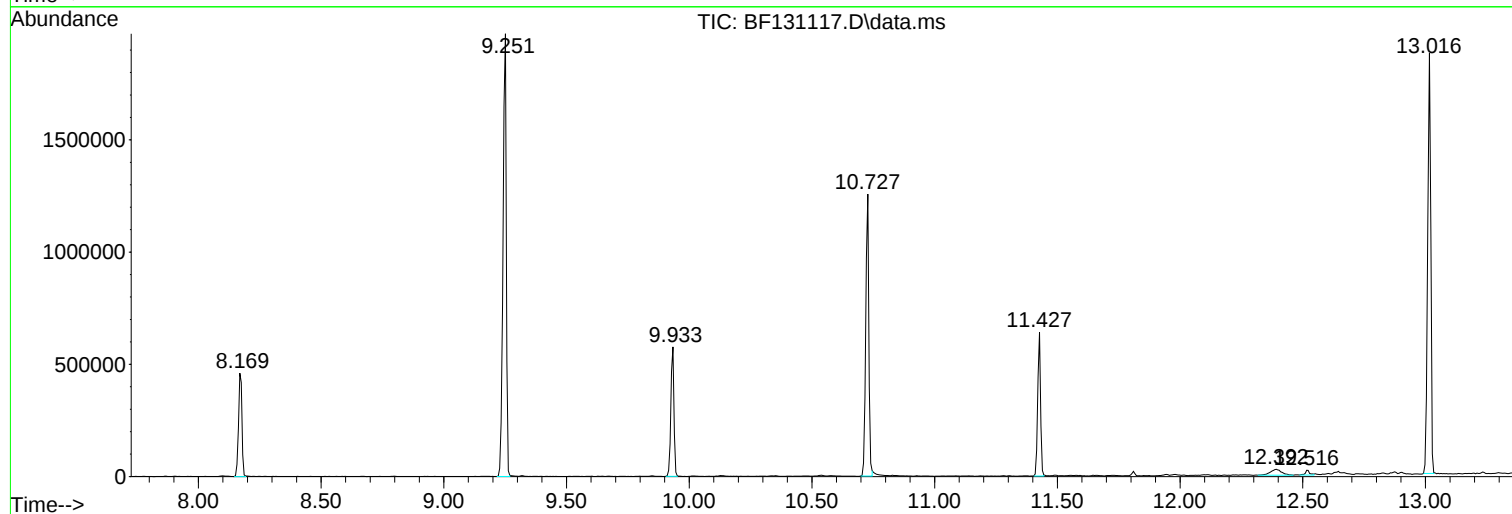
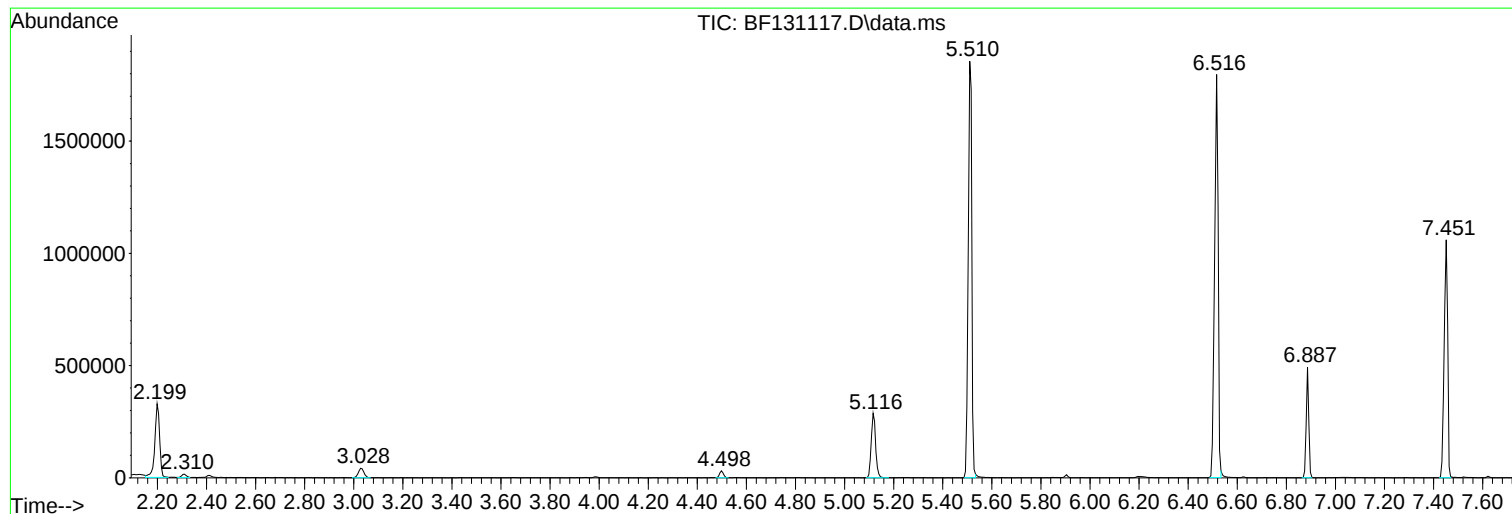
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF110922.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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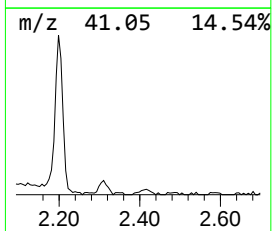
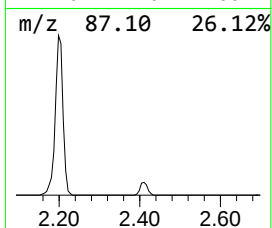
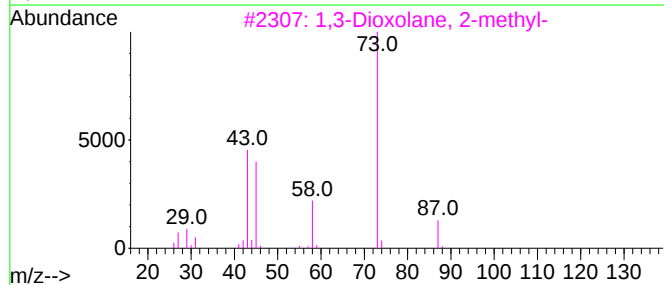
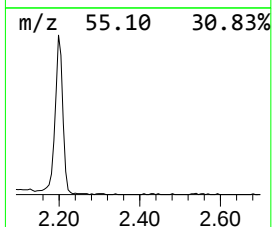
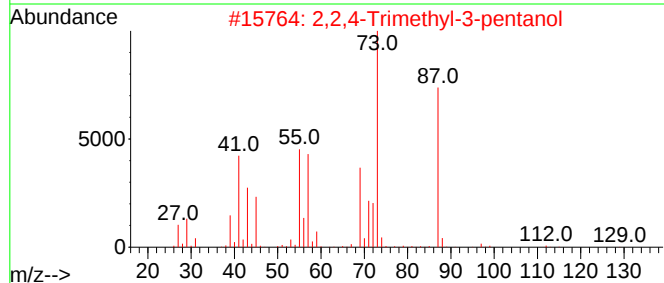
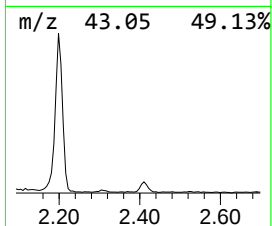
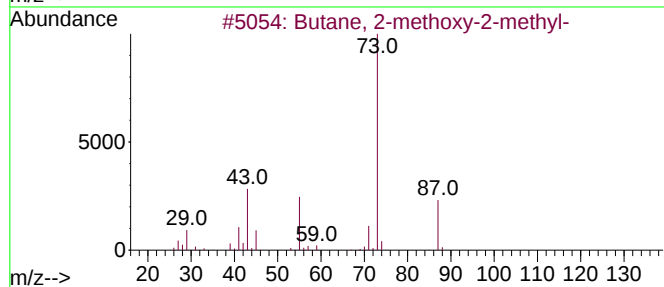
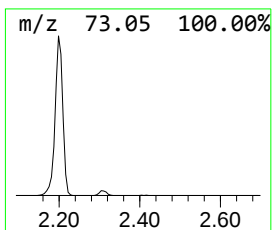
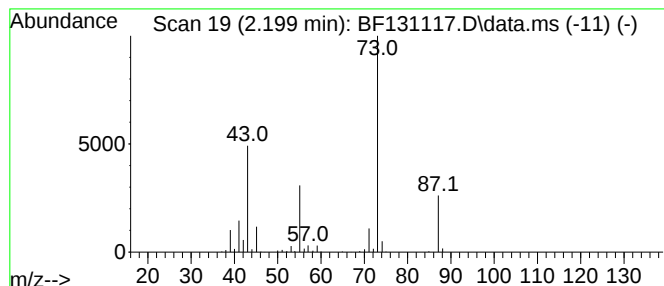
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.199	24.92 ng	442270	1,4-Dichlorobenzene-d4	6.887

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			1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	37
4			Octanal, 7-methoxy-3,7-dimethyl-	186	C11H22O2	003613-30-7	17
5			Silane, tetramethyl-	88	C4H12Si	000075-76-3	9



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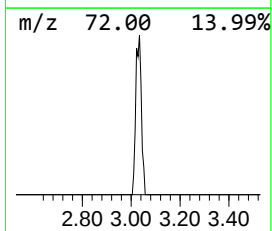
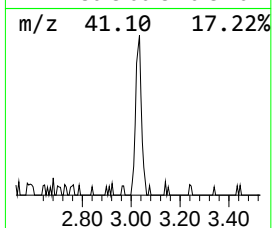
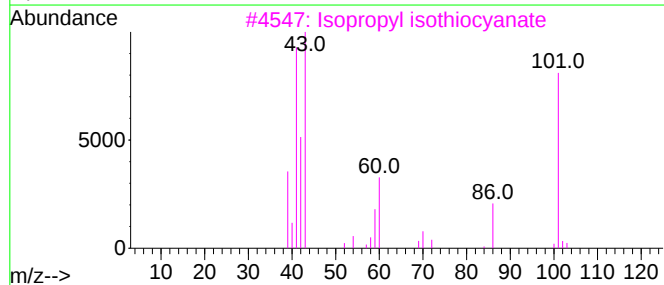
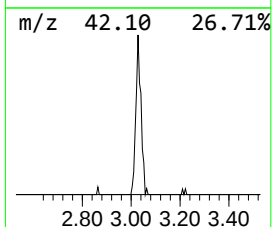
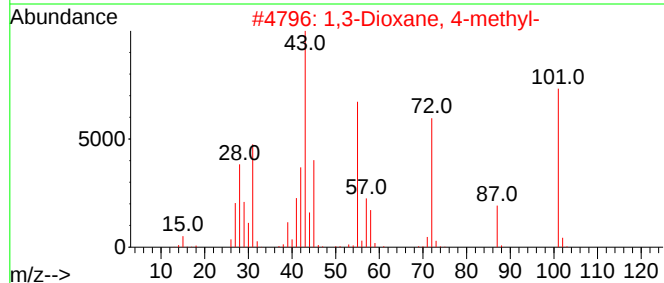
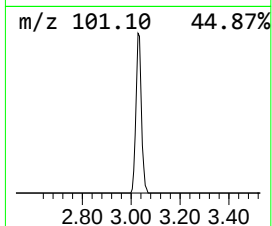
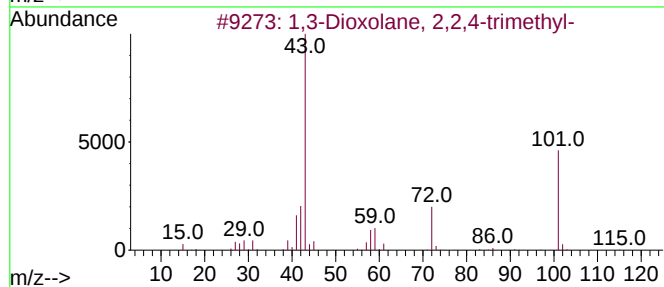
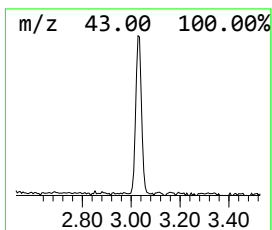
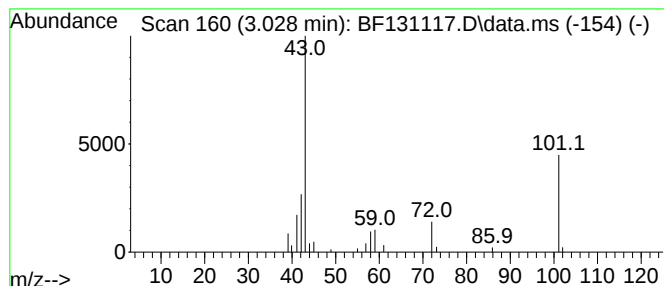
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TIC Library : C:\Database\NIST20.L
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Peak Number 2 1,3-Dioxolane, 2,2,4-trimet... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.028	3.75 ng	66514	1,4-Dichlorobenzene-d4	6.887

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,3-Dioxolane, 2,2,4-trimethyl-	116	C6H12O2	001193-11-9	72	
2	1,3-Dioxane, 4-methyl-	102	C5H10O2	001120-97-4	42	
3	Isopropyl isothiocyanate	101	C4H7NS	002253-73-8	36	
4	1-Methoxy-2-propyl acetate	132	C6H12O3	000108-65-6	10	
5	Butane	58	C4H10	000106-97-8	9	



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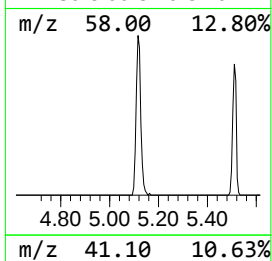
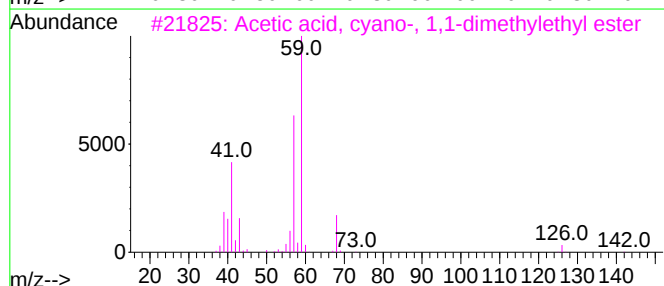
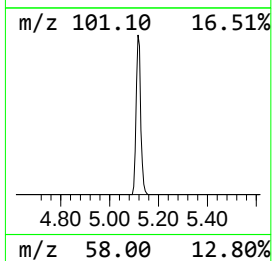
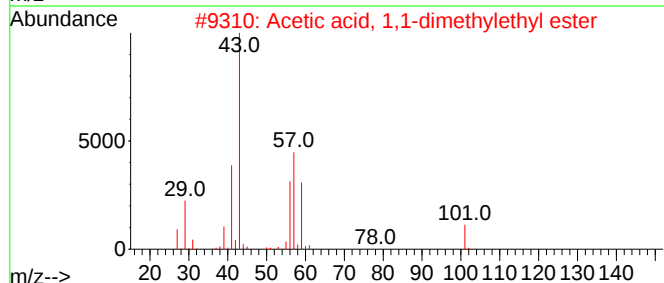
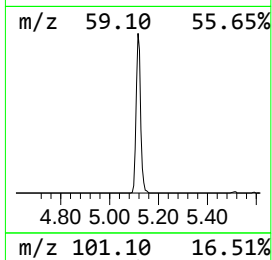
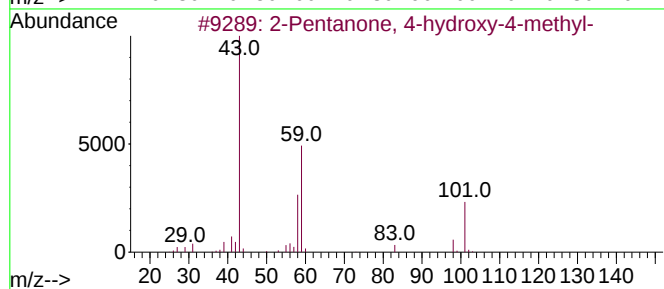
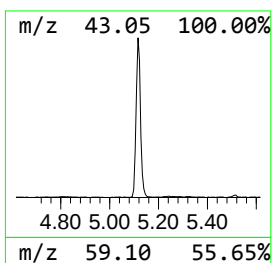
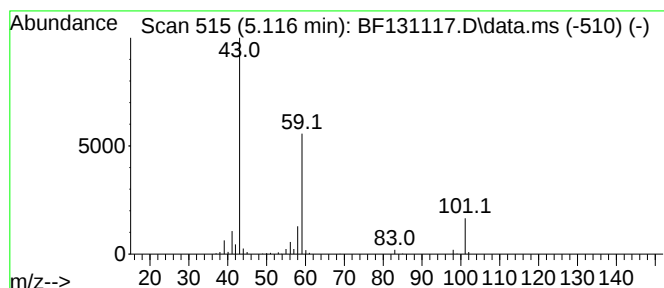
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.116	20.17 ng	357856	1,4-Dichlorobenzene-d4	6.887

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	72
2			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	32
3			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	32
4			1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	12
5			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9



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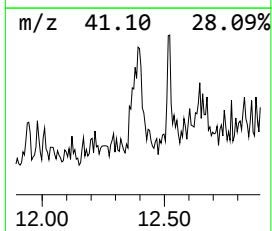
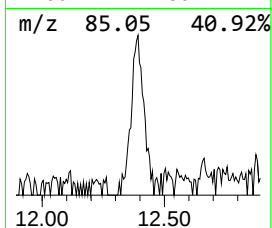
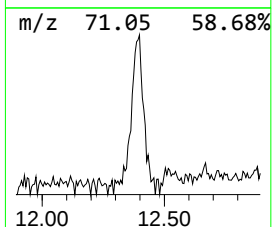
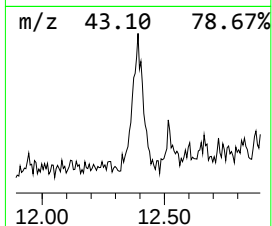
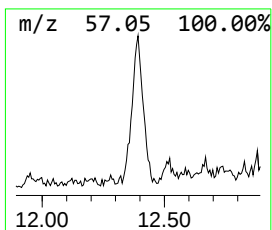
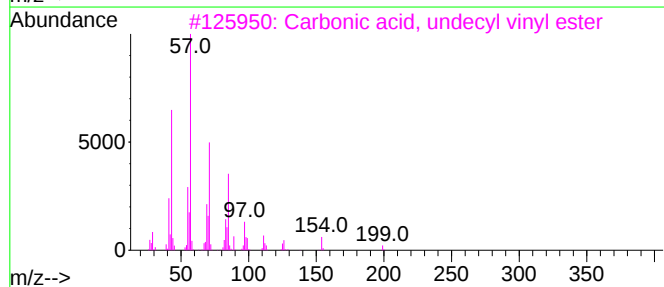
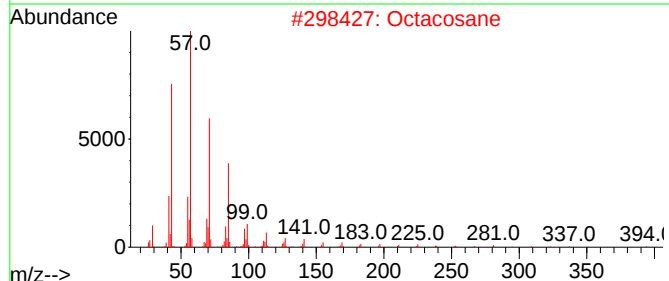
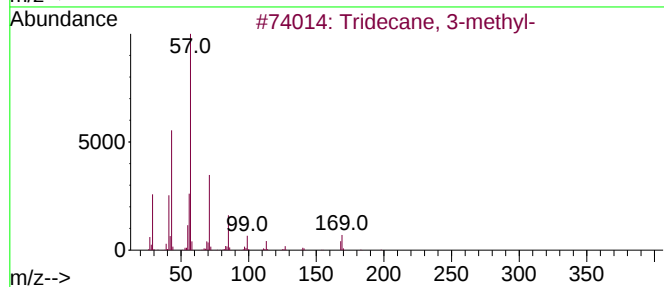
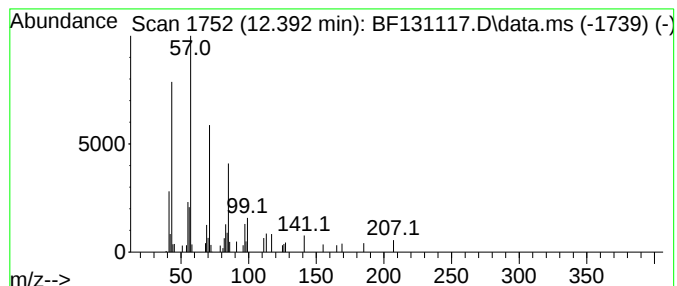
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Peak Number 4 Tridecane, 3-methyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.392	3.37 ng	86721	Phenanthrene-d10	11.427

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tridecane, 3-methyl-	198	C14H30	006418-41-3	80
2			Octacosane	394	C28H58	000630-02-4	80
3			Carbonic acid, undecyl vinyl ester	242	C14H26O3	1000382-54-9	80
4			Hentriacontane	437	C31H64	000630-04-6	80
5			10-Methylnonadecane	282	C20H42	056862-62-5	74



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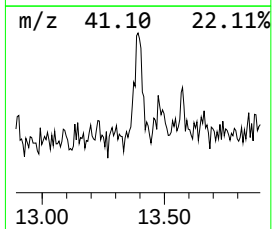
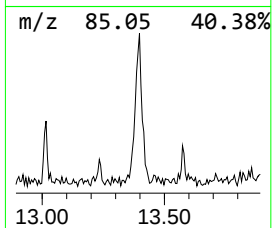
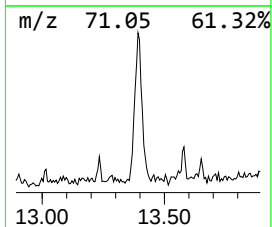
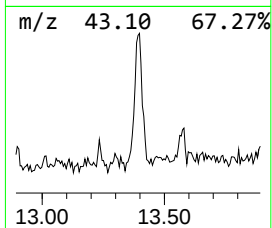
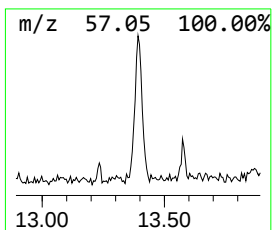
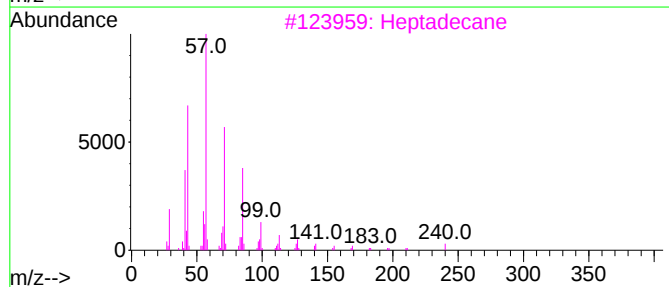
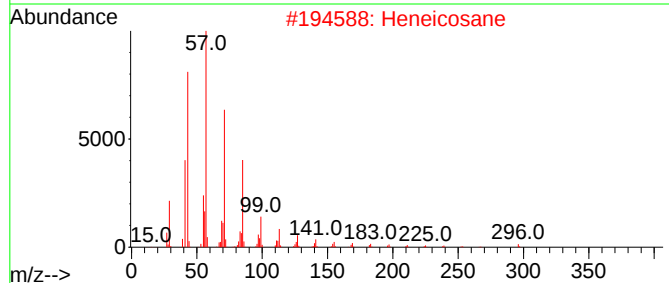
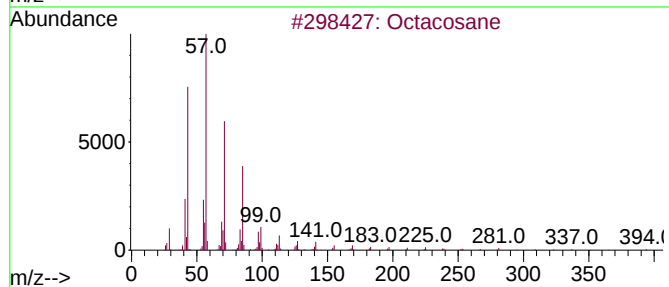
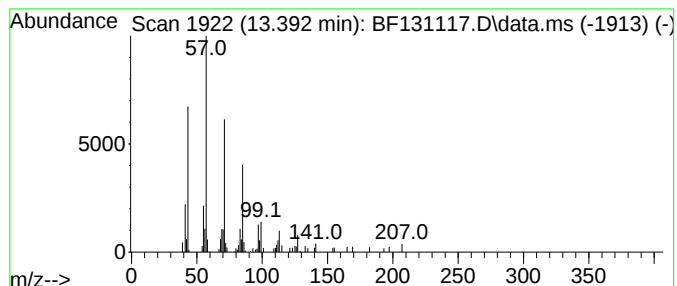
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Peak Number 5 Octacosane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.392	4.81 ng	101734	Chrysene-d12	14.074

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octacosane	394	C28H58	000630-02-4	90
2			Heneicosane	296	C21H44	000629-94-7	87
3			Heptadecane	240	C17H36	000629-78-7	87
4			Tetracosane	338	C24H50	000646-31-1	87
5			Heptacosane	380	C27H56	000593-49-7	86



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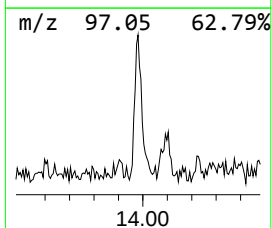
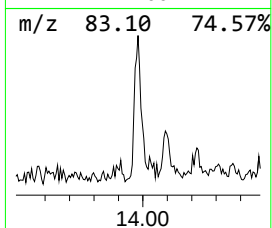
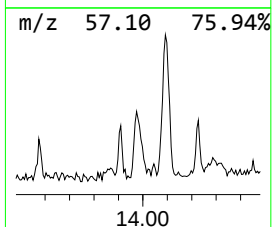
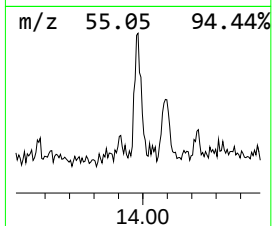
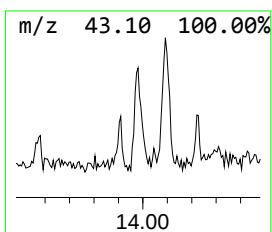
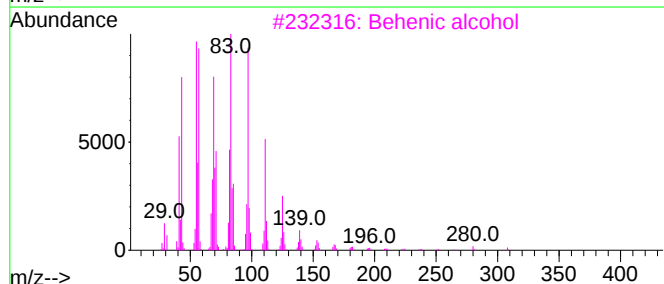
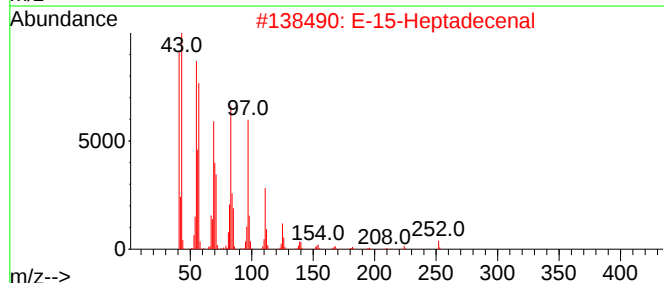
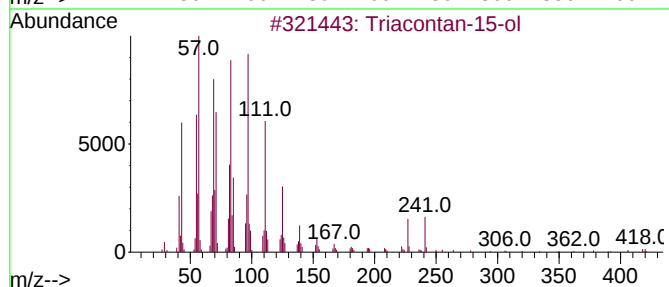
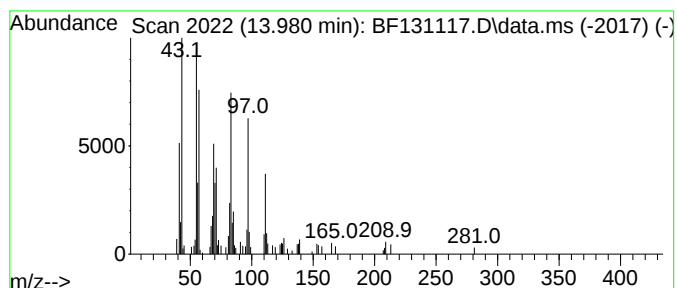
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Triacontan-15-ol Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.980	4.57 ng	96834	Chrysene-d12	14.074

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Triacontan-15-ol	438	C30H62O	031849-26-0	91
2			E-15-Heptadecenal	252	C17H32O	1000130-97-9	83
3			Behenic alcohol	326	C22H46O	000661-19-8	83
4			1-Hexacosene	364	C26H52	018835-33-1	81
5			1-Tetracosene	336	C24H48	010192-32-2	81



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF111022\
Data File : BF131117.D
Acq On : 10 Nov 2022 11:51
Operator : CG\JU
Sample : N5495-10
Misc :
ALS Vial : 4 Sample Multiplier: 1

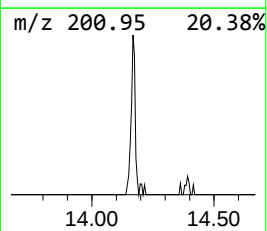
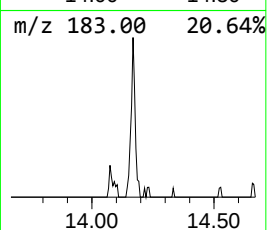
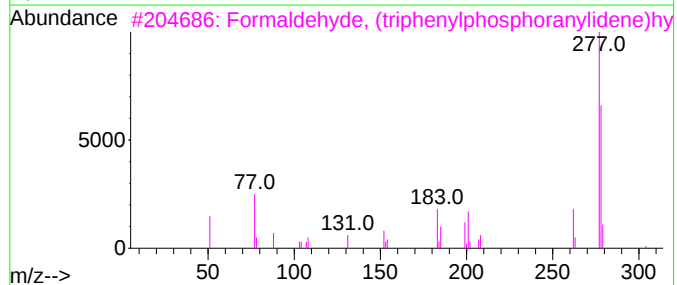
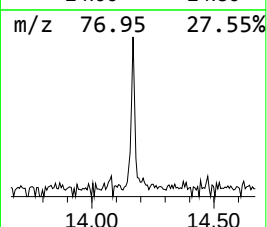
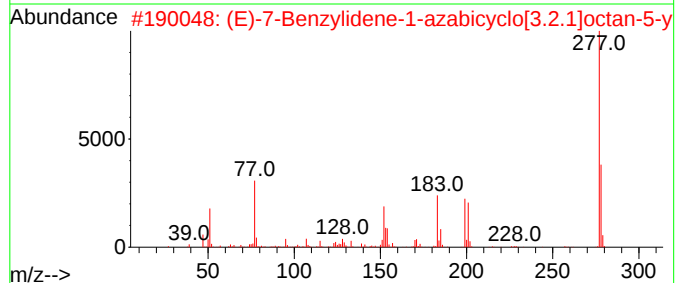
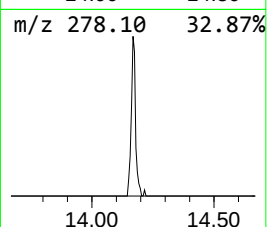
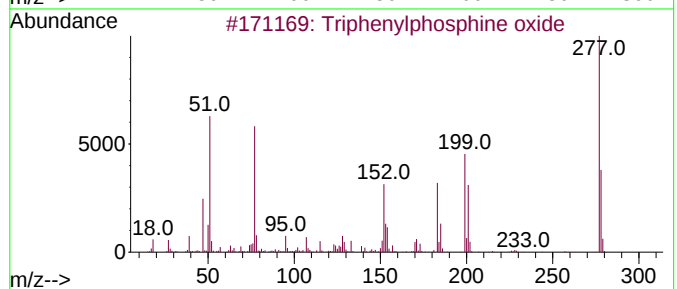
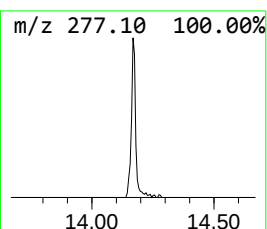
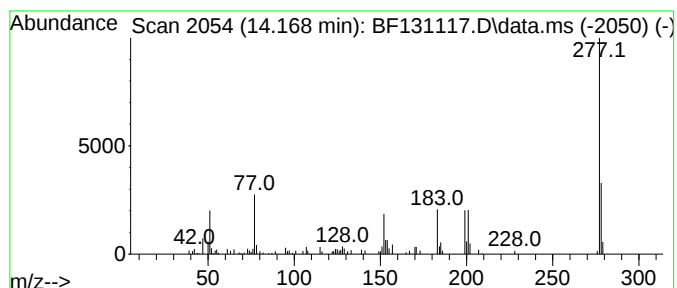
Instrument :
BNA_F
ClientSampleId :
DUP-01

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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 Triphenylphosphine oxide Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.168	2.49 ng	52765	Chrysene-d12	14.074	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Triphenylphosphine oxide	278	C18H15OP	000791-28-6	96
2	(E)-7-Benzylidene-1-azabicyclo[3...	293	C15H19NO3S	1000483-83-4	64
3	Formaldehyde, (triphenylphosphor...	304	C19H17N2P	015990-54-2	39
4	2-Phenyl-6-(trifluoromethyl)-1H-...	277	C15H10F3NO	1000387-59-3	38
5	1H-Indole-3-acetic acid, 5-[(tri...	277	C14H19NO3Si	072101-35-0	27



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF111022\
Data File : BF131117.D
Acq On : 10 Nov 2022 11:51
Operator : CG\JU
Sample : N5495-10
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
DUP-01

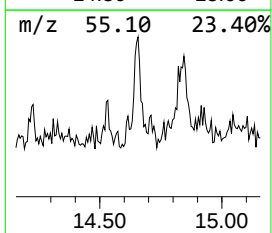
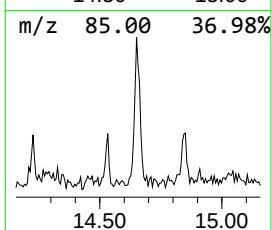
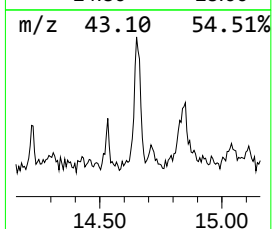
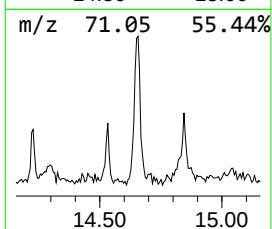
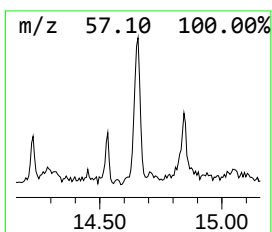
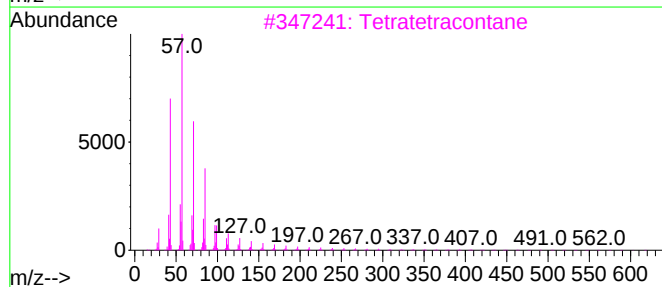
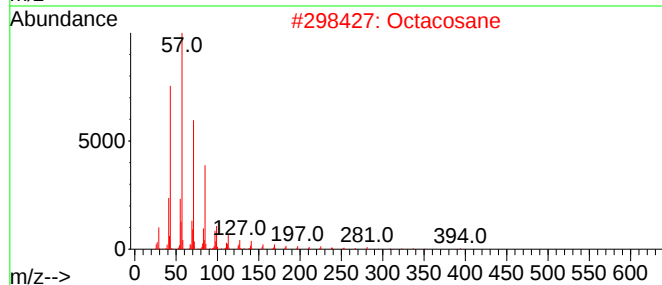
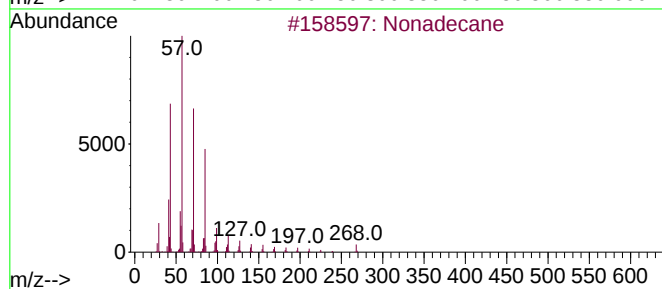
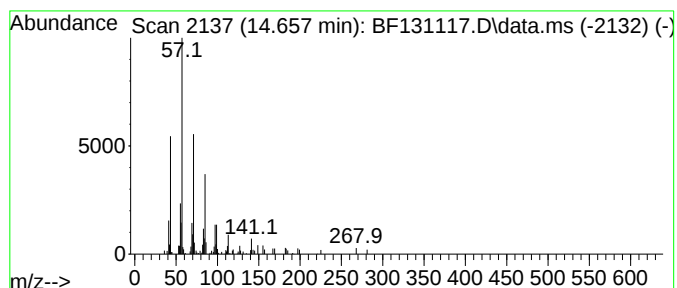
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Nonadecane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.657	3.58 ng	75809	Chrysene-d12	14.074

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonadecane	268	C19H40	000629-92-5	96
2			Octacosane	394	C28H58	000630-02-4	90
3			Tetratetracontane	619	C44H90	007098-22-8	87
4			Octacosane, 2-methyl-	408	C29H60	001560-98-1	87
5			Hentriacontane	437	C31H64	000630-04-6	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF111022\
Data File : BF131117.D
Acq On : 10 Nov 2022 11:51
Operator : CG\JU
Sample : N5495-10
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
DUP-01

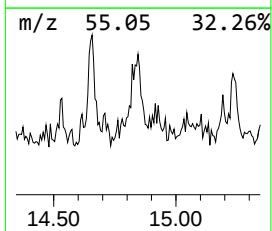
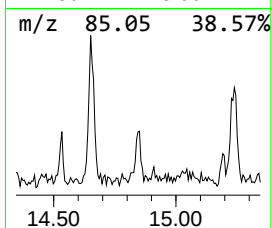
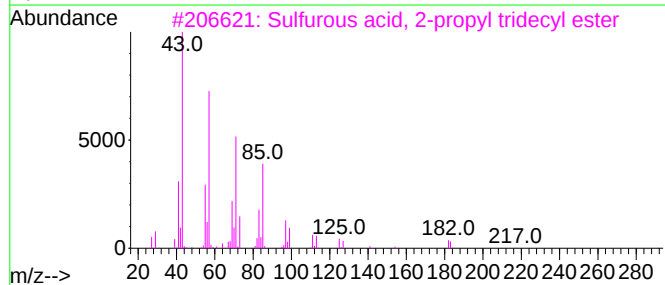
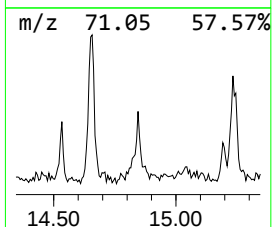
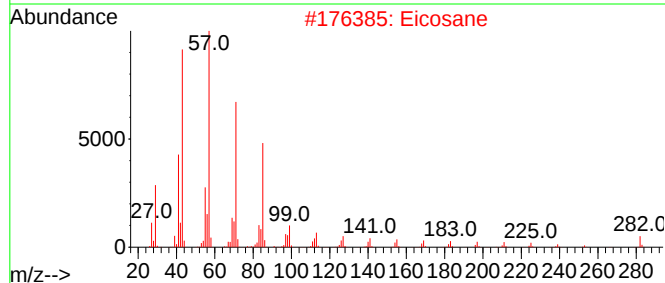
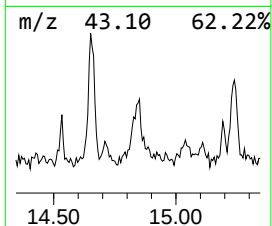
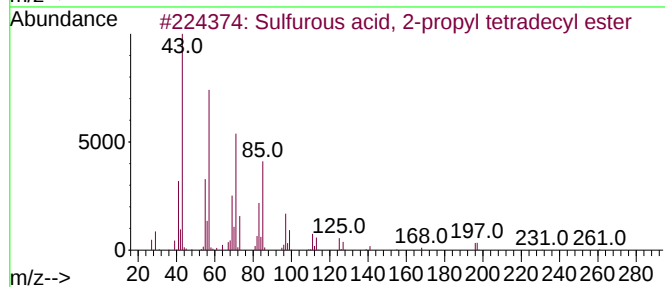
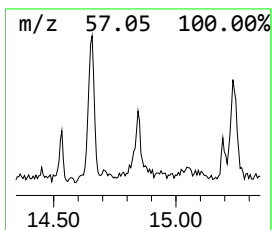
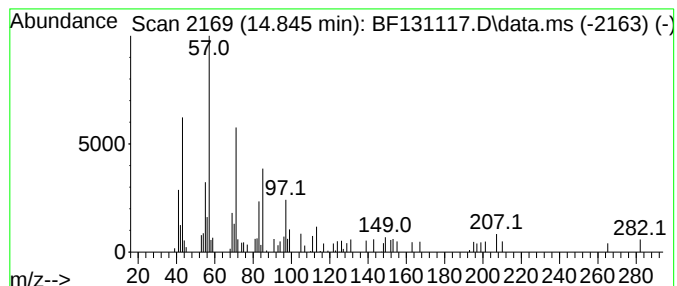
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Sulfurous acid, 2-propyl te... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.845	3.28 ng	53817	Perylene-d12	15.574

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Sulfurous acid, 2-propyl tetradecyl ester	320	C17H36O3S	1010309-12-5	80
2			Eicosane	282	C20H42	000112-95-8	76
3			Sulfurous acid, 2-propyl tridecyl ester	306	C16H34O3S	1000309-12-4	72
4			Carbonic acid, octadecyl vinyl ester	340	C21H40O3	1000382-54-4	72
5			Sulfurous acid, butyl hexadecyl ester	362	C20H42O3S	1000309-18-3	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF111022\
Data File : BF131117.D
Acq On : 10 Nov 2022 11:51
Operator : CG\JU
Sample : N5495-10
Misc :
ALS Vial : 4 Sample Multiplier: 1

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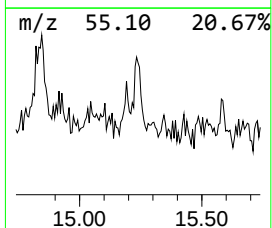
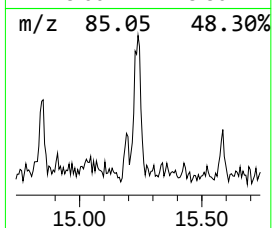
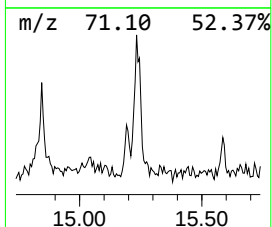
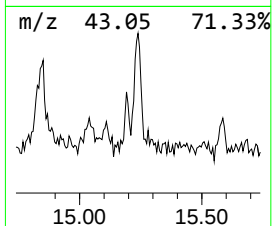
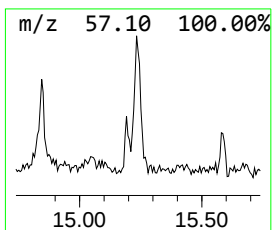
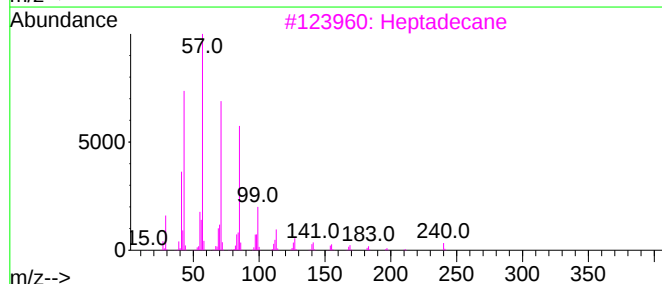
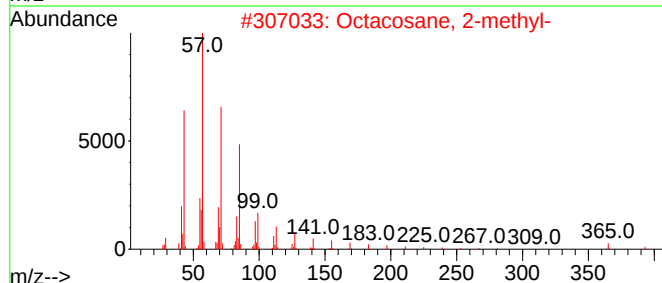
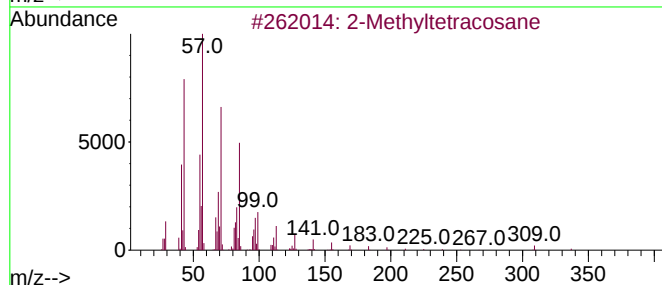
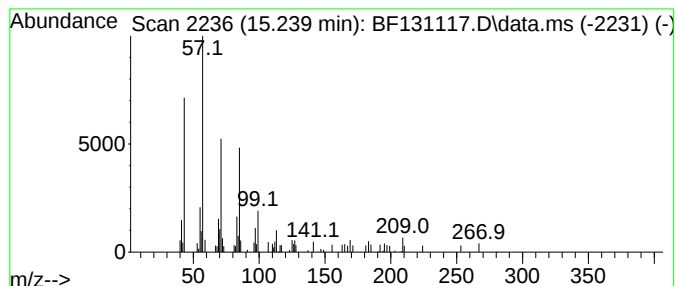
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 2-Methyltetracosane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.239	3.47 ng	56921	Perylene-d12	15.574

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		2-Methyltetracosane	352	C25H52	001560-78-7	86
2		Octacosane, 2-methyl-	408	C29H60	001560-98-1	86
3		Heptadecane	240	C17H36	000629-78-7	86
4		Octacosane	394	C28H58	000630-02-4	74
5		Hexatriacontane	507	C36H74	000630-06-8	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF111022\
Data File : BF131117.D
Acq On : 10 Nov 2022 11:51
Operator : CG\JU
Sample : N5495-10
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
DUP-01

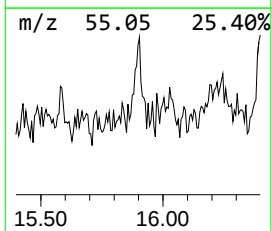
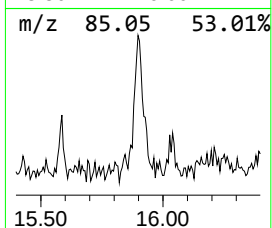
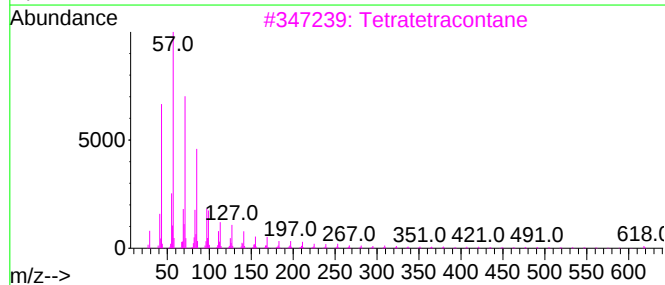
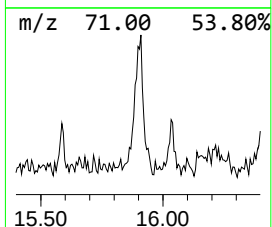
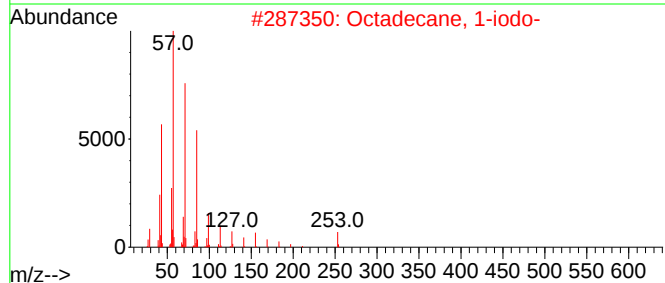
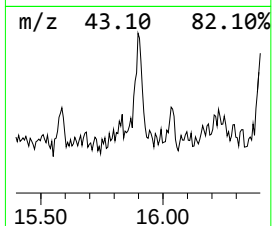
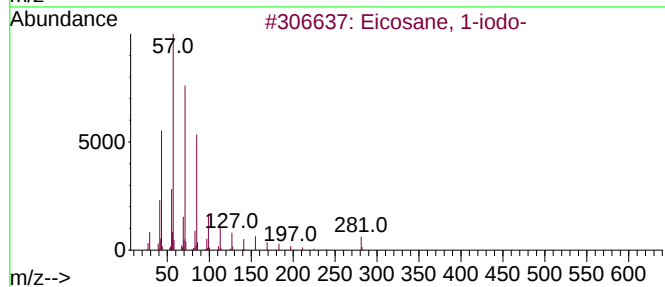
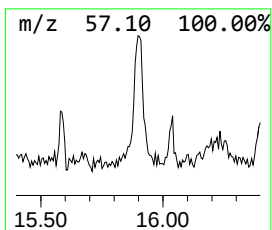
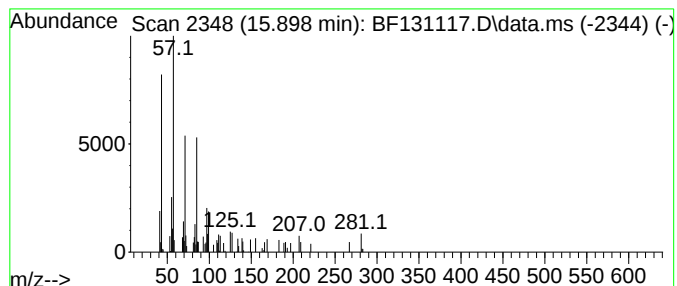
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Eicosane, 1-iodo- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.898	3.57 ng	58640	Perylene-d12	15.574

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Eicosane, 1-iodo-	408	C20H41I	1000406-31-8	59
2		Octadecane, 1-iodo-	380	C18H37I	000629-93-6	53
3		Tetratetracontane	619	C44H90	007098-22-8	50
4		Eicosane	282	C20H42	000112-95-8	46
5		Hexadecane, 2,6,10,14-tetramethyl-	282	C20H42	000638-36-8	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF111022\
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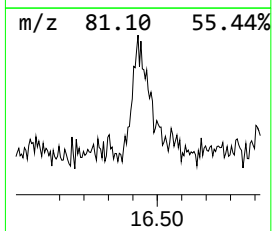
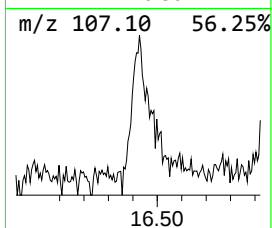
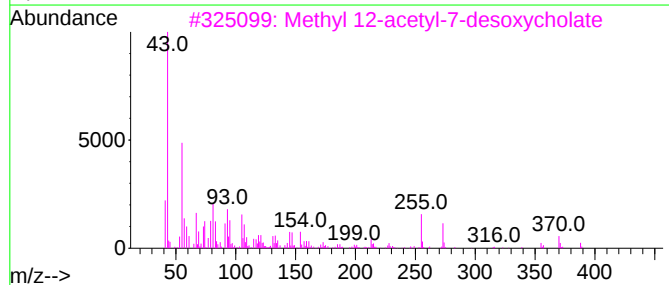
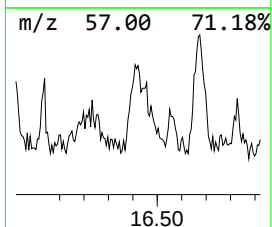
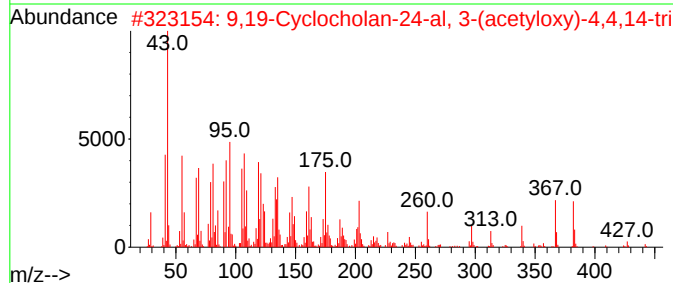
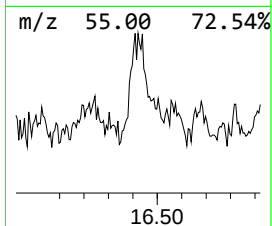
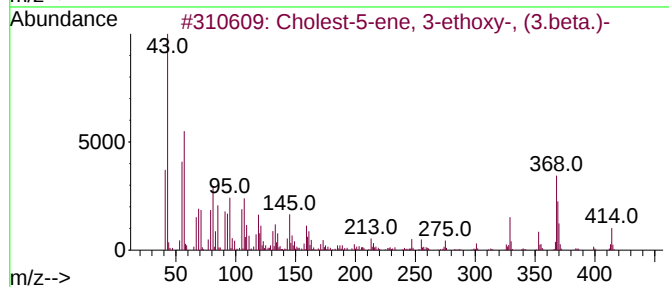
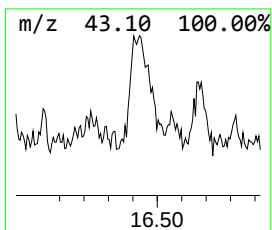
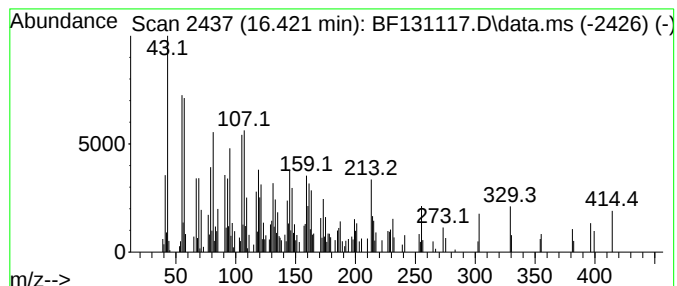
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF110922.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Cholest-5-ene, 3-ethoxy-, (... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.		
16.421	23.42 ng	384493	Perylene-d12	15.574		
Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cholest-5-ene, 3-ethoxy-, (3.bet...	414	C29H50O	000986-19-6	58
2		9,19-Cyclochohan-24-al, 3-(acety...	442	C29H46O3	025089-90-1	37
3		Methyl 12-acetyl-7-desoxycholate	448	C27H44O5	055547-48-3	25
4		Stigmastan-3-en-6-ol	414	C29H50O	1000214-19-7	16
5		Stigmastane, 23,24-epoxy-, (5.al...	414	C29H50O	054411-91-5	16



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF111022\
Data File : BF131117.D
Acq On : 10 Nov 2022 11:51
Operator : CG\JU
Sample : N5495-10
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
DUP-01

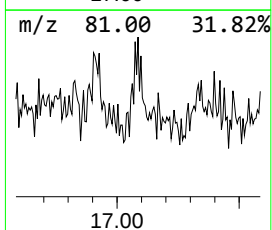
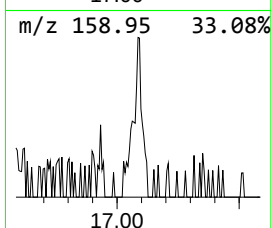
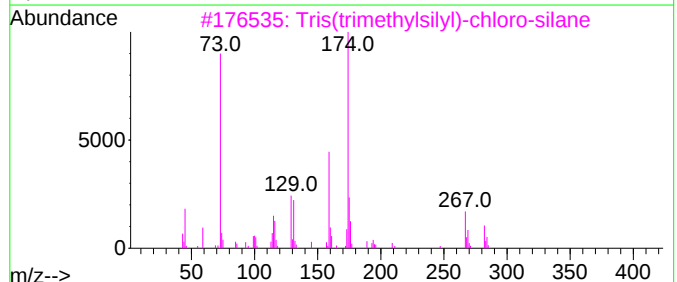
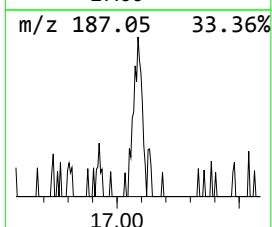
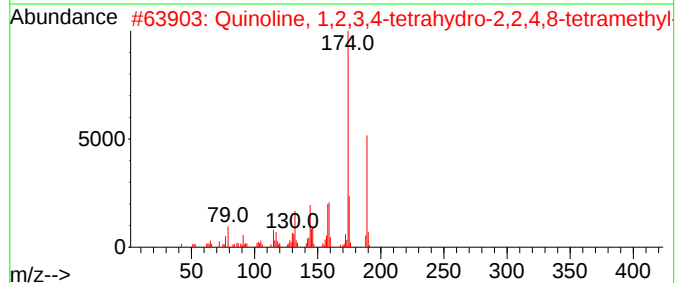
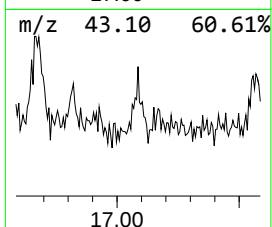
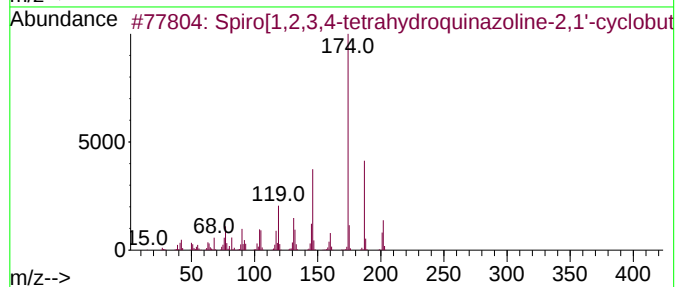
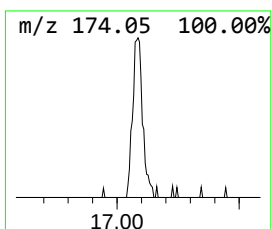
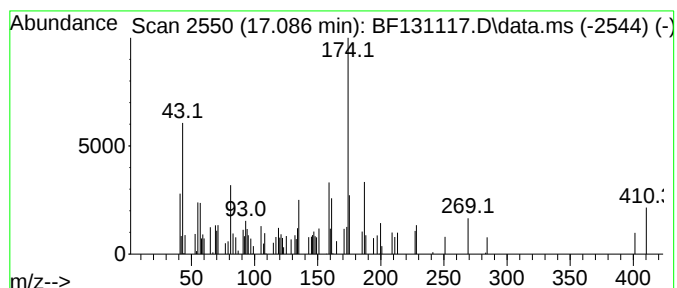
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 unknown17.086 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.086	3.25 ng	53278	Perylene-d12	15.574

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Spiro[1,2,3,4-tetrahydroquinazol...	202	C12H14N2O	1000508-42-3	38
2			Quinoline, 1,2,3,4-tetrahydro-2,...	189	C13H19N	1000308-78-8	38
3			Tris(trimethylsilyl)-chloro-silane	282	C9H27ClSi4	005565-32-2	37
4			2,8,9-Trioxa-5-aza-1-silabicyclo...	203	C8H17N3Si	002097-16-7	35
5			Tyramine, 3TMS derivative	353	C17H35NOSi3	068595-84-6	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF111022\
Data File : BF131117.D
Acq On : 10 Nov 2022 11:51
Operator : CG\JU
Sample : N5495-10
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
DUP-01

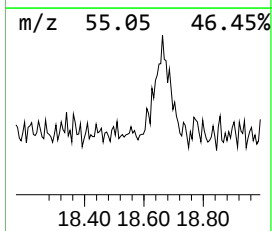
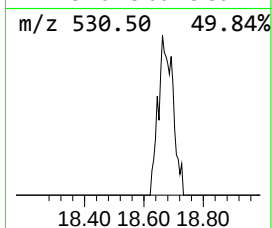
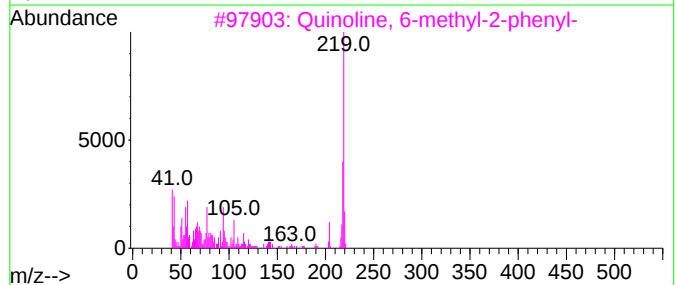
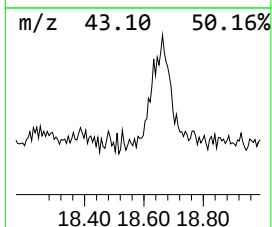
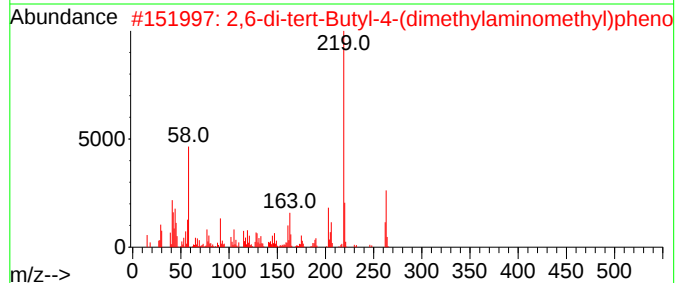
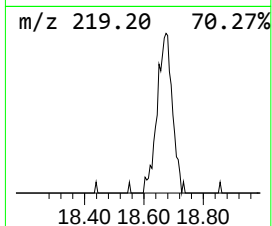
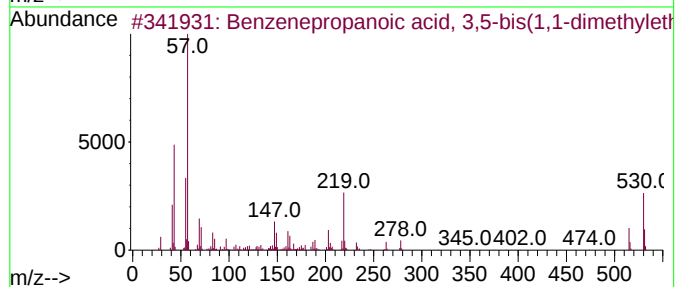
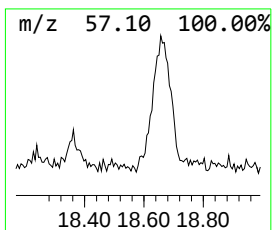
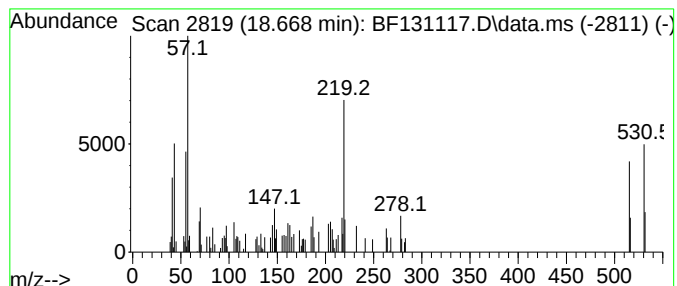
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Benzenepropanoic acid, 3,5-... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.668	6.28 ng	103157	Perylene-d12	15.574

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzenepropanoic acid, 3,5-bis(1...	530	C35H62O3	002082-79-3	53
2			2,6-di-tert-Butyl-4-(dimethylami...	263	C17H29NO	000088-27-7	18
3			Quinoline, 6-methyl-2-phenyl-	219	C16H13N	027356-46-3	11
4			5-(3,4-Dimethoxyphenyl)-2H-pyraz...	219	C11H13N3O2	1000409-59-6	11
5			3,5-di-tert-Butyl-4-hydroxybenza...	234	C15H22O2	001620-98-0	10



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF111022\
Data File : BF131117.D
Acq On : 10 Nov 2022 11:51
Operator : CG\JU
Sample : N5495-10
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
DUP-01

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF110922.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...	2.199	24.9	ng	442270	1	6.887	354910	20.0
1,3-Dioxolane, ...	3.028	3.8	ng	66514	1	6.887	354910	20.0
2-Pentanone, 4-...	5.116	20.2	ng	357856	1	6.887	354910	20.0
Tridecane, 3-me...	12.392	3.4	ng	86721	4	11.427	515044	20.0
Octacosane	13.392	4.8	ng	101734	5	14.074	423381	20.0
Triacontan-15-ol	13.980	4.6	ng	96834	5	14.074	423381	20.0
Triphenylphosph...	14.168	2.5	ng	52765	5	14.074	423381	20.0
Nonadecane	14.657	3.6	ng	75809	5	14.074	423381	20.0
Sulfurous acid,...	14.845	3.3	ng	53817	6	15.574	328318	20.0
2-Methyltetraaco...	15.239	3.5	ng	56921	6	15.574	328318	20.0
Eicosane, 1-iodo-	15.898	3.6	ng	58640	6	15.574	328318	20.0
Cholest-5-ene, ...	16.421	23.4	ng	384493	6	15.574	328318	20.0
unknown17.086	17.086	3.3	ng	53278	6	15.574	328318	20.0
Benzenepropanoi...	18.668	6.3	ng	103157	6	15.574	328318	20.0