

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF122922\
 Data File : BF131877.D
 Acq On : 29 Dec 2022 14:44
 Operator : CG\JU
 Sample : N6232-08
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MW-2

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-BF122422.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF131877.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.110	338	344	349	rBV	25912	33619	1.04%	0.184%
2	5.016	494	498	504	rBV	255845	277831	8.60%	1.520%
3	5.428	564	568	579	rBV	1021117	983407	30.44%	5.379%
4	6.445	737	741	750	rBV	719780	615768	19.06%	3.368%
5	6.816	800	804	808	rBV	632127	545297	16.88%	2.982%
6	7.392	891	902	904	rBV	1492308	1831937	56.71%	10.020%
7	7.563	925	931	934	rBV2	26601	35422	1.10%	0.194%
8	7.698	936	954	957	rBV	381304	1114145	34.49%	6.094%
9	7.851	975	980	986	rVB2	89809	88294	2.73%	0.483%
10	8.034	1005	1011	1015	rBV6	36054	69906	2.16%	0.382%
11	8.104	1016	1023	1026	rBV	716235	728943	22.57%	3.987%
12	8.257	1045	1049	1054	rVB3	56768	64003	1.98%	0.350%
13	8.328	1054	1061	1064	rBV3	87515	123904	3.84%	0.678%
14	8.381	1068	1070	1077	rVB7	25661	46594	1.44%	0.255%
15	8.445	1077	1081	1083	rBV3	31132	39853	1.23%	0.218%
16	8.492	1083	1089	1093	rBV6	35828	70340	2.18%	0.385%
17	8.587	1099	1105	1107	rVB2	38719	50049	1.55%	0.274%
18	8.651	1113	1116	1117	rBV2	44158	40221	1.25%	0.220%
19	8.681	1117	1121	1125	rVV3	53867	93815	2.90%	0.513%
20	8.828	1142	1146	1147	rBV3	34678	44030	1.36%	0.241%
21	8.863	1150	1152	1159	rVB4	60346	85564	2.65%	0.468%
22	8.969	1166	1170	1173	rBV4	43080	63879	1.98%	0.349%
23	9.004	1173	1176	1180	rVB	51510	52144	1.61%	0.285%
24	9.192	1199	1208	1210	rBV	3407549	3230273	100.00%	17.668%
25	9.216	1210	1212	1217	rVV4	54491	50631	1.57%	0.277%
26	9.304	1222	1227	1232	rVB6	86270	112677	3.49%	0.616%
27	9.392	1238	1242	1247	rVB2	135601	152448	4.72%	0.834%
28	9.457	1252	1253	1258	rVB4	33464	37835	1.17%	0.207%
29	9.498	1258	1260	1263	rBV3	30717	41366	1.28%	0.226%
30	9.534	1263	1266	1269	rVB3	49433	54095	1.67%	0.296%
31	9.728	1296	1299	1303	rBV4	27109	39584	1.23%	0.216%
32	9.839	1315	1318	1319	rBV2	41572	44926	1.39%	0.246%
33	9.869	1319	1323	1326	rVV	956393	803359	24.87%	4.394%
34	10.086	1358	1360	1364	rVB2	54879	53729	1.66%	0.294%
35	10.151	1369	1371	1375	rBV3	37406	43542	1.35%	0.238%
36	10.339	1400	1403	1405	rBV4	36178	43051	1.33%	0.235%

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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-BF122422.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

37	10.369	1406	1408	1411	rVV3	39410	40160	1.24%	0.220%
38	10.498	1428	1430	1432	rBV3	38767	40828	1.26%	0.223%
39	10.669	1453	1459	1462	rBV	1869593	1849339	57.25%	10.115%
40	11.186	1545	1547	1549	rVB2	52898	38331	1.19%	0.210%
41	11.363	1573	1577	1580	rBV	939932	775469	24.01%	4.241%
42	11.751	1641	1643	1646	rVB3	47272	34949	1.08%	0.191%
43	11.898	1665	1668	1675	rVB	193992	196855	6.09%	1.077%
44	12.680	1799	1801	1808	rVB2	46767	50634	1.57%	0.277%
45	12.957	1844	1848	1851	rBV	2045733	1875250	58.05%	10.256%
46	13.857	1998	2001	2009	rBV	198913	218274	6.76%	1.194%
47	14.016	2024	2028	2033	rVB	590846	508107	15.73%	2.779%
48	14.104	2040	2043	2045	rBV	83728	91826	2.84%	0.502%
49	14.127	2045	2047	2051	rVB	266567	241453	7.47%	1.321%
50	15.498	2275	2280	2291	rVB	350436	467784	14.48%	2.558%
51	16.015	2365	2368	2376	rVB8	23463	47906	1.48%	0.262%

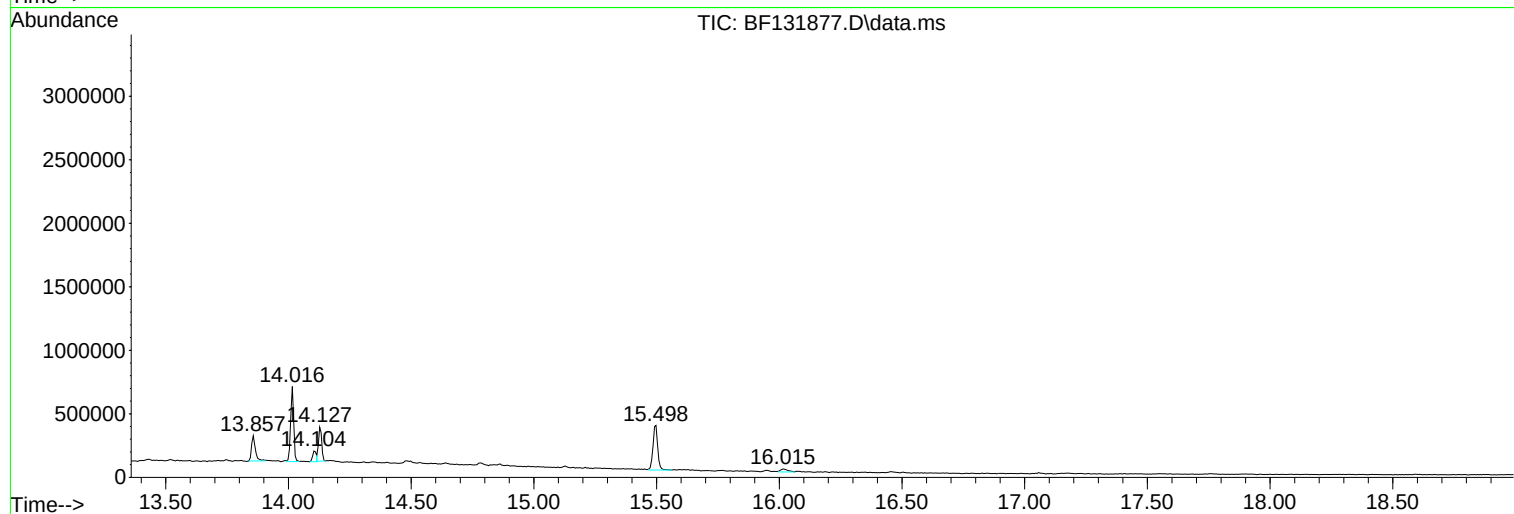
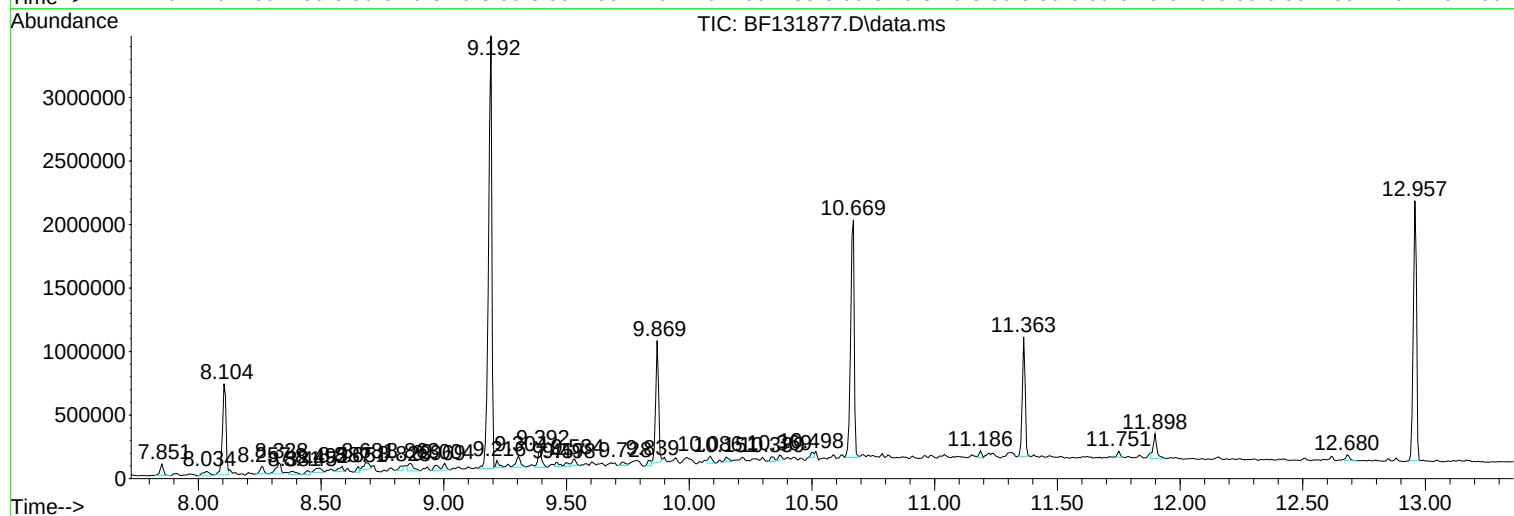
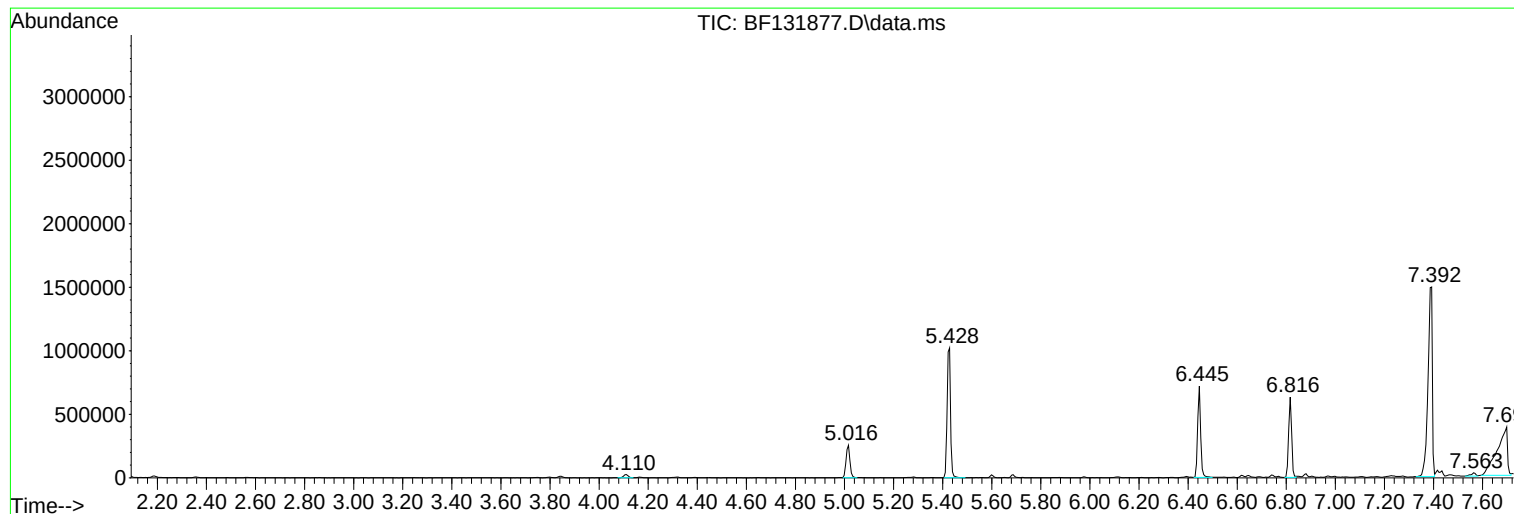
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF122422.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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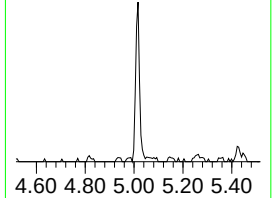
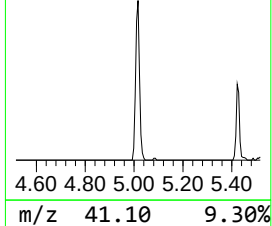
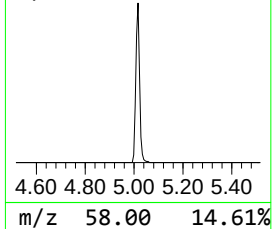
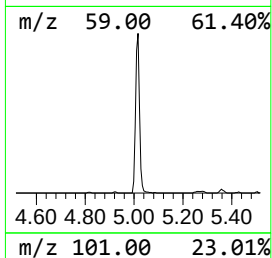
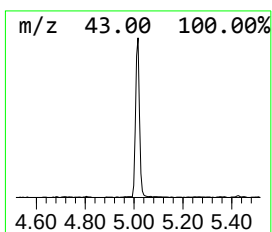
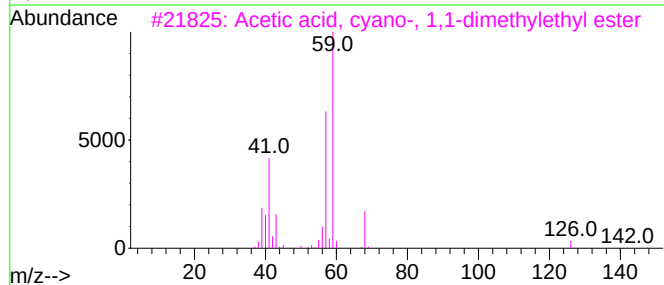
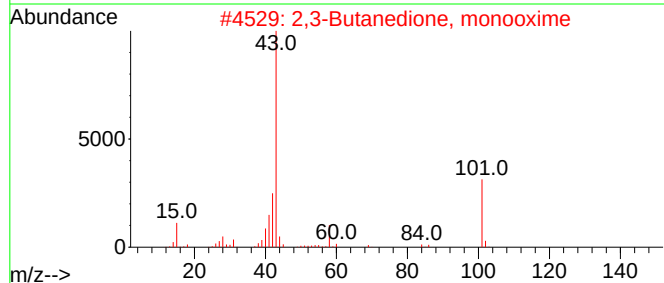
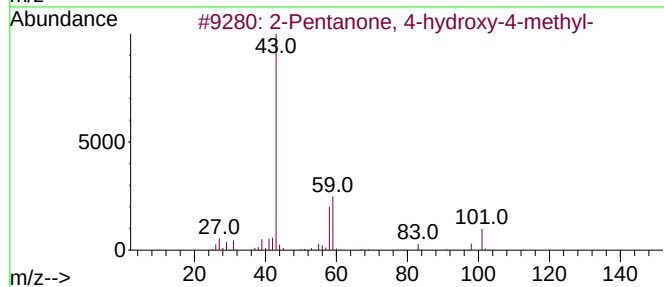
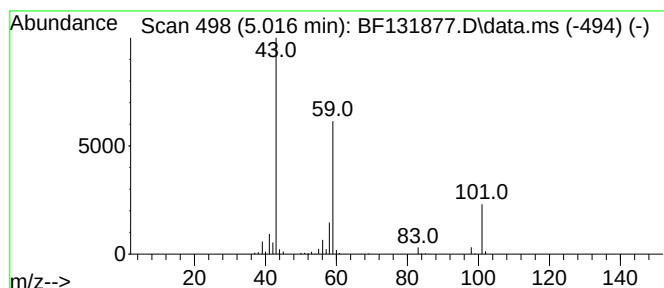
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF122422.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.016	10.19 ng	277831	1,4-Dichlorobenzene-d4	6.816

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	72		
2	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	25		
3	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	23		
4	2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	23		
5	1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10		



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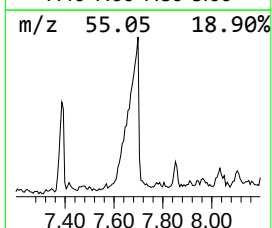
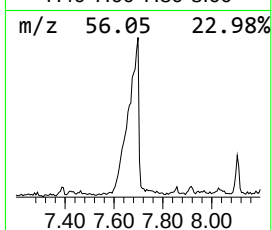
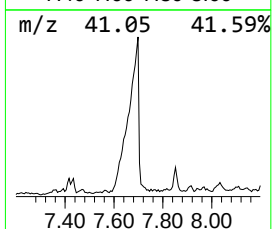
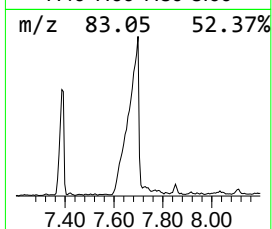
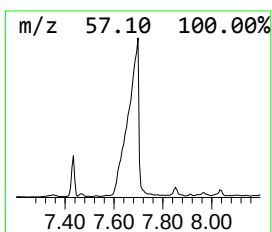
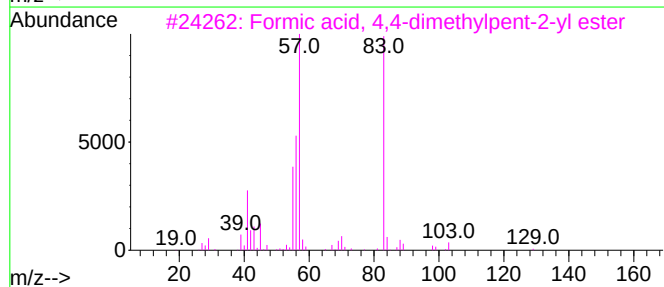
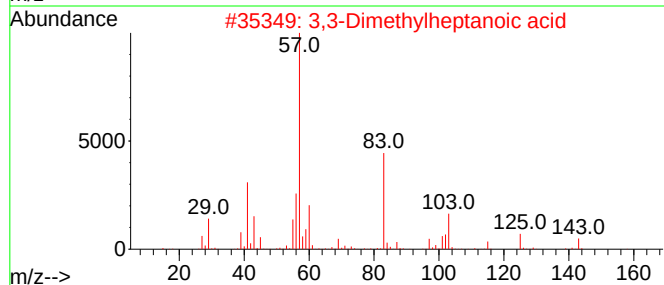
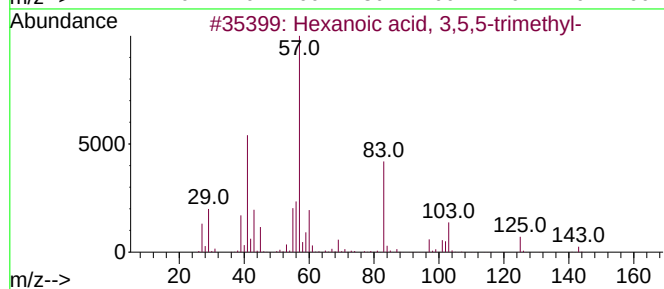
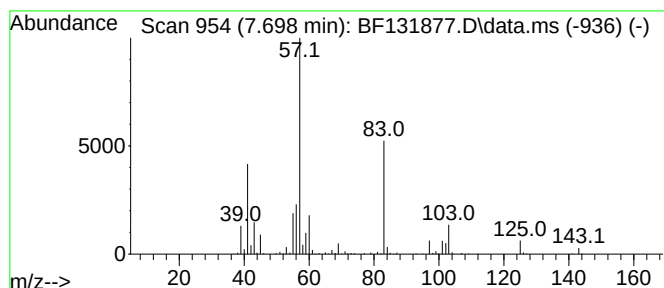
Instrument :
BNA_F
ClientSampleId :
MW-2

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF122422.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 Hexanoic acid, 3,5,5-trimet... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.	
7.698	30.57 ng	1114150	Naphthalene-d8	8.104	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexanoic acid, 3,5,5-trimethyl-	158	C9H18O2	003302-10-1	80
2	3,3-Dimethylheptanoic acid	158	C9H18O2	067061-30-7	50
3	Formic acid, 4,4-dimethylpent-2-...	144	C8H16O2	1000368-69-9	43
4	Hexanoyl chloride, 3,5,5-trimethyl-	176	C9H17ClO	036727-29-4	42
5	4,4-Dimethyl-2-pentanol, chlorod...	228	C9H15ClF2O2	1000376-27-0	32



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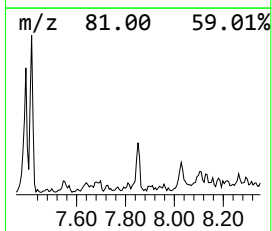
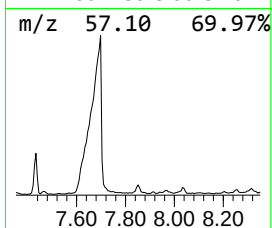
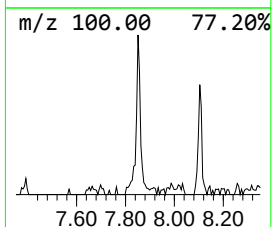
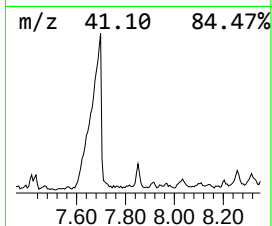
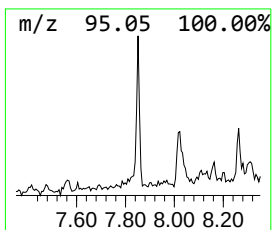
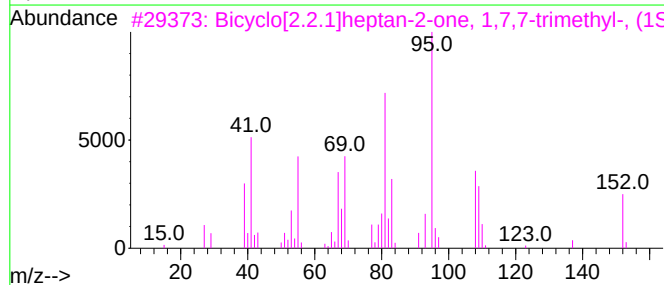
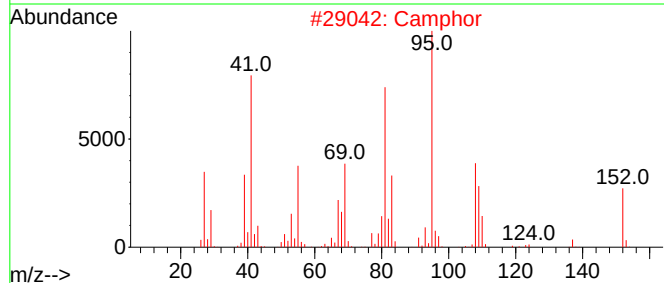
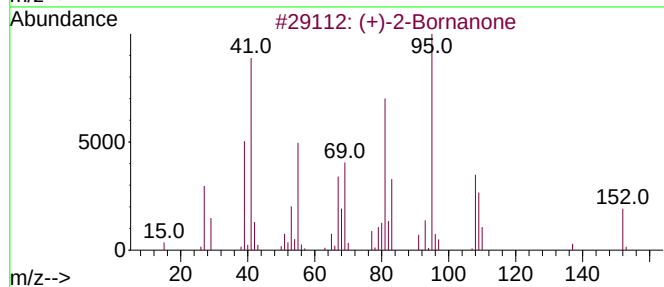
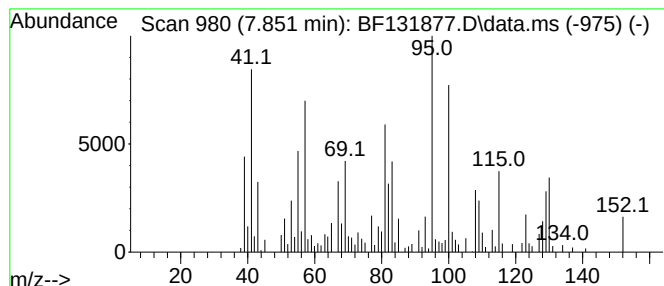
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 (+)-2-Bornanone Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.851	2.42 ng	88294	Naphthalene-d8	8.104

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			(+)-2-Bornanone	152	C10H16O	000464-49-3	86
2			Camphor	152	C10H16O	000076-22-2	38
3			Bicyclo[2.2.1]heptan-2-one, 1,7,...	152	C10H16O	000464-48-2	35
4			5-Heptenoic acid, 2,6-dimethyl-2...	182	C11H18O2	100052-75-3	27
5			Ethyl 2-methylcyclopropanecarbox...	128	C7H12O2	020913-25-1	27



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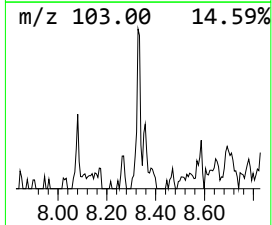
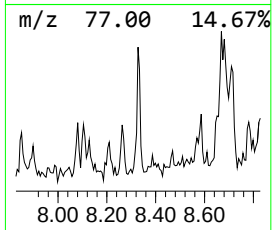
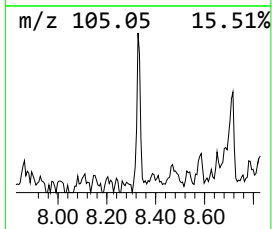
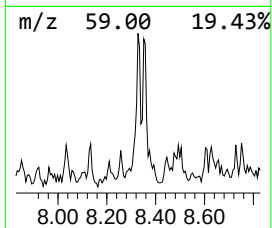
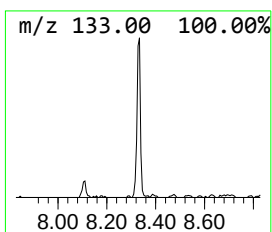
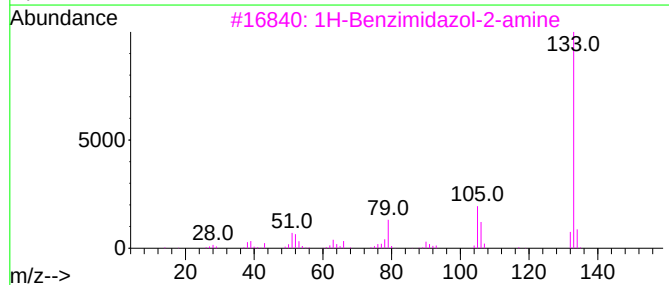
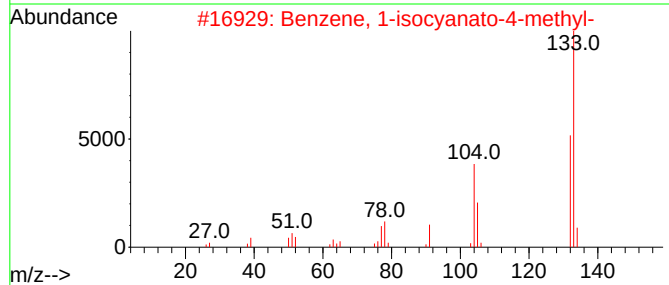
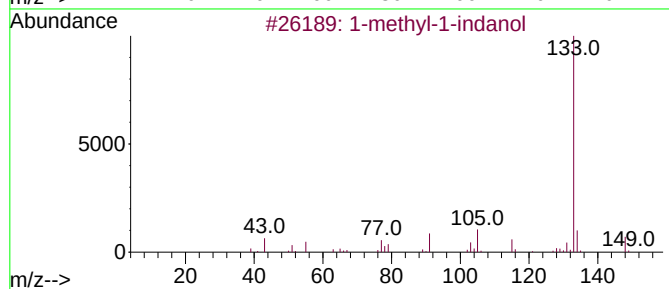
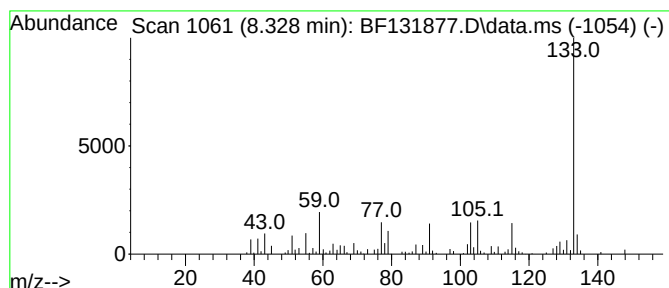
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 4 1-methyl-1-indanol Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.328	3.40 ng	123904	Naphthalene-d8	8.104

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		1-methyl-1-indanol	148	C10H12O	064666-42-8	76
2		Benzene, 1-isocyanato-4-methyl-	133	C8H7NO	000622-58-2	60
3		1H-Benzimidazol-2-amine	133	C7H7N3	000934-32-7	53
4		Propan-1-one, 3-(4-chlorophenyl)...	414	C23H20Cl2OS	284493-90-9	52
5		Benzene, 1,3-dimethyl-5-(1-methy...	148	C11H16	004706-90-5	50



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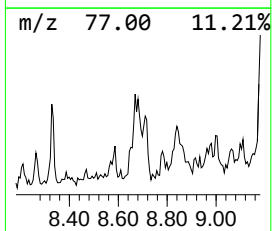
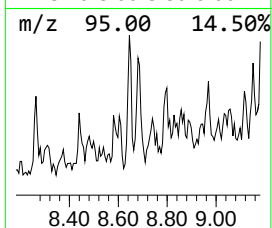
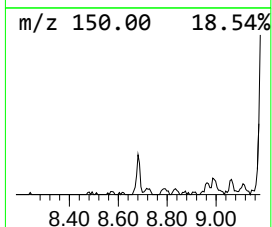
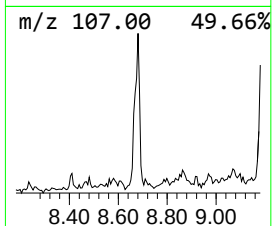
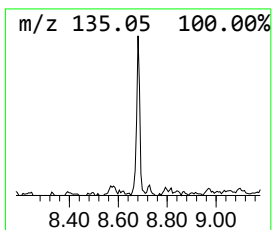
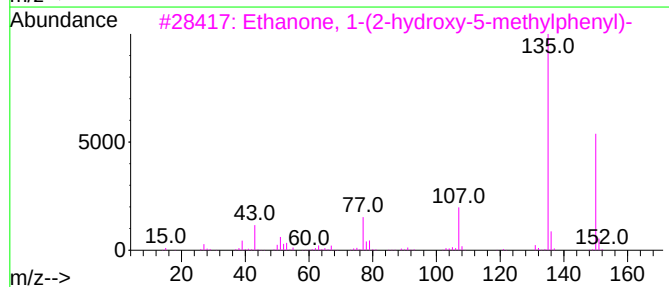
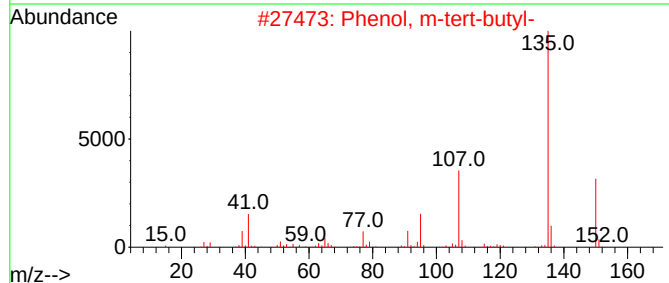
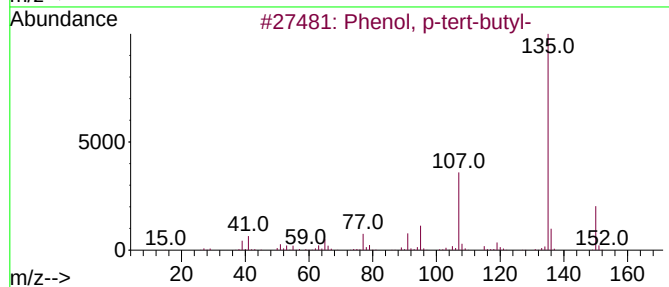
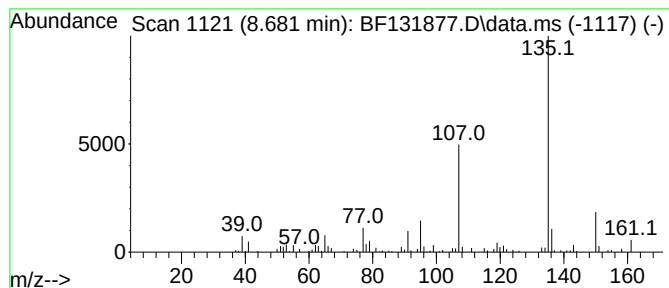
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Phenol, p-tert-butyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.681	2.57 ng	93815	Naphthalene-d8	8.104

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	91
2			Phenol, m-tert-butyl-	150	C10H14O	000585-34-2	91
3			Ethanone, 1-(2-hydroxy-5-methylp...	150	C9H10O2	001450-72-2	58
4			Imidazo[1,2-c]pyrimidin-5-ol	135	C6H5N3O	055662-66-3	58
5			Cyclohexene, 2-ethenyl-1,3,3-tri...	150	C11H18	005293-90-3	58



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF122922\
Data File : BF131877.D
Acq On : 29 Dec 2022 14:44
Operator : CG\JU
Sample : N6232-08
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MW-2

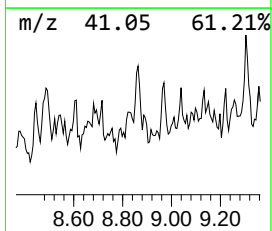
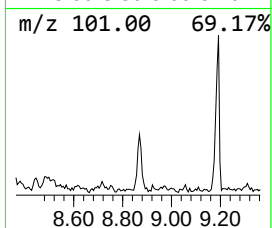
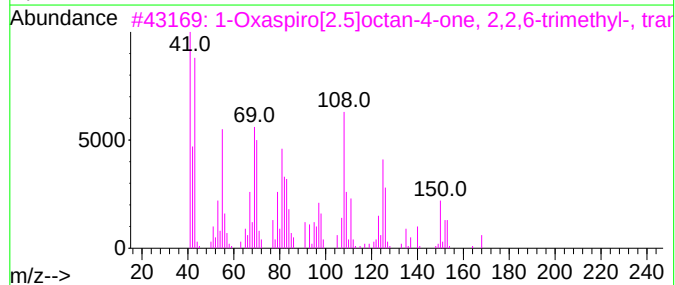
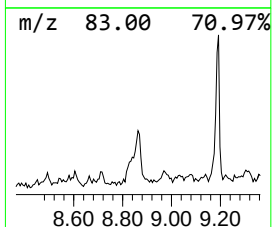
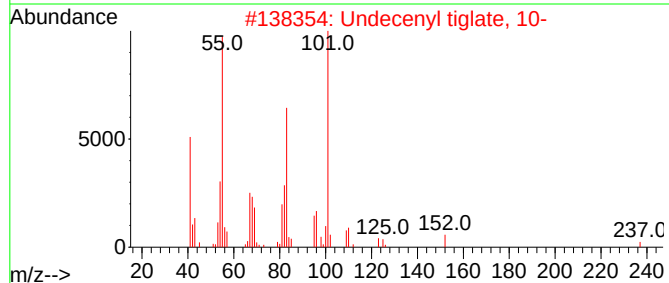
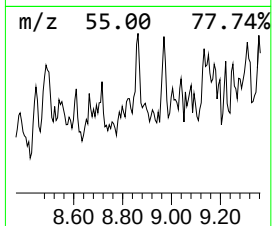
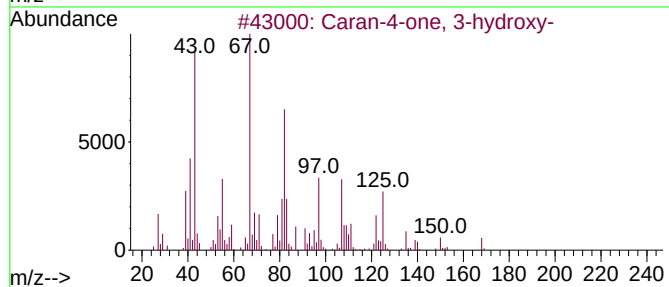
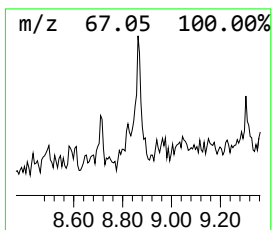
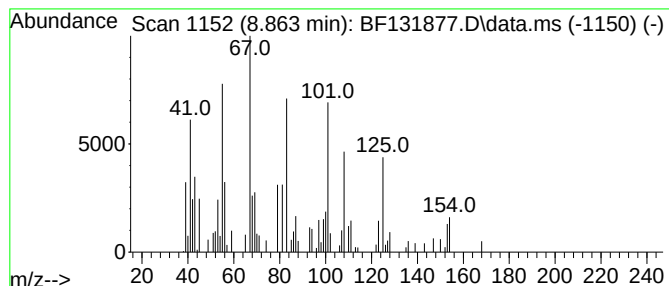
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 unknown8.863 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.863	2.35 ng	85564	Naphthalene-d8	8.104

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Caran-4-one, 3-hydroxy-	168	C10H16O2	024348-06-9	35
2			Undecenyl tiglate, 10-	252	C16H28O2	1000383-56-4	22
3			1-Oxaspiro[2.5]octan-4-one, 2,2,...	168	C10H16O2	013080-28-9	20
4			Decan-2-yl (E)-2-methylbut-2-enoate	240	C15H28O2	1000373-74-7	14
5			4-Chloro-3-n-hexyltetrahydropyran	204	C11H21ClO	066555-66-6	12



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF122922\
Data File : BF131877.D
Acq On : 29 Dec 2022 14:44
Operator : CG\JU
Sample : N6232-08
Misc :
ALS Vial : 5 Sample Multiplier: 1

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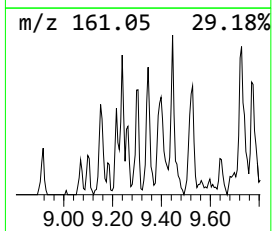
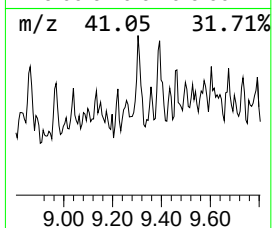
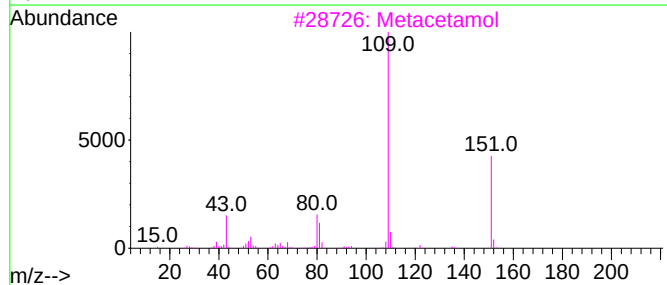
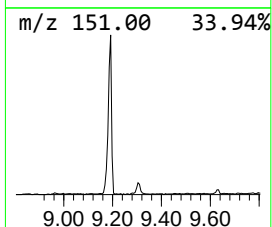
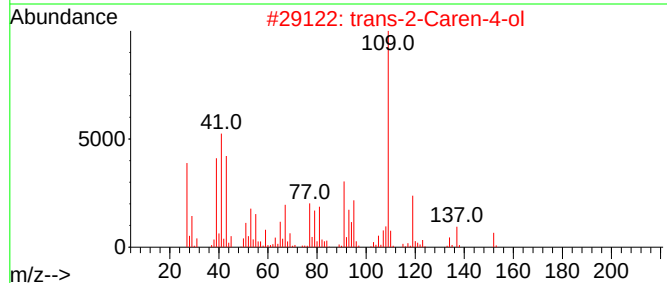
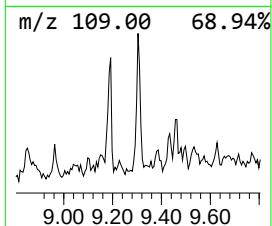
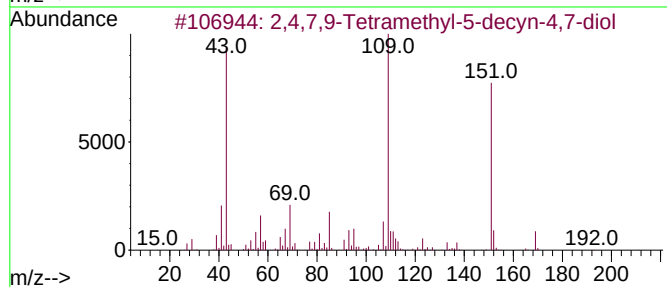
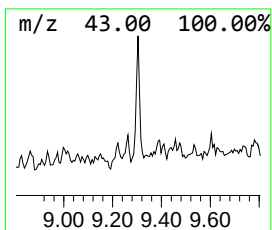
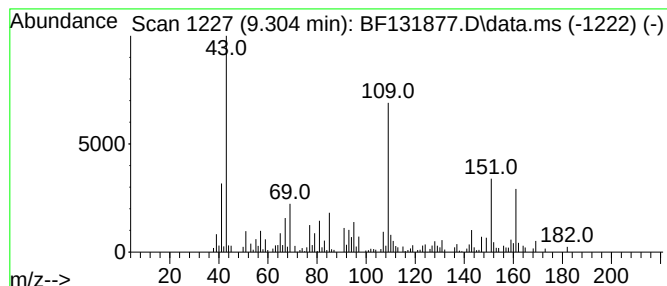
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 unknown9.304 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.304	2.81 ng	112677	Acenaphthene-d10	9.869

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,4,7,9-Tetramethyl-5-decyn-4,7-...	226	C14H26O2	000126-86-3	47		
2	trans-2-Caren-4-ol	152	C10H16O	004017-82-7	46		
3	Metacetamol	151	C8H9NO2	000621-42-1	46		
4	Benzenamine, 4-propoxy-	151	C9H13NO	004469-80-1	35		
5	Pyridine, 3-butyl-, 1-oxide	151	C9H13NO	031396-33-5	35		



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF122922\
Data File : BF131877.D
Acq On : 29 Dec 2022 14:44
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Sample : N6232-08
Misc :
ALS Vial : 5 Sample Multiplier: 1

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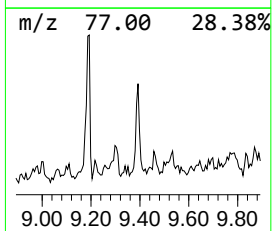
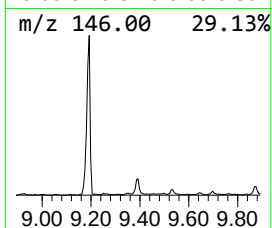
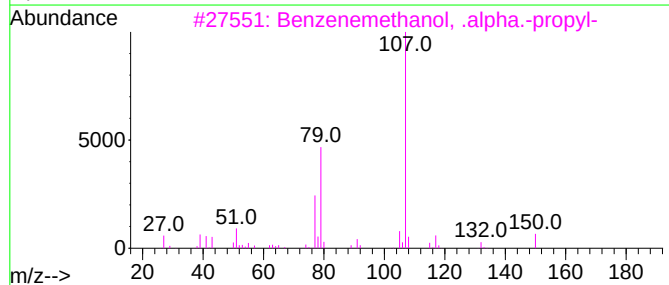
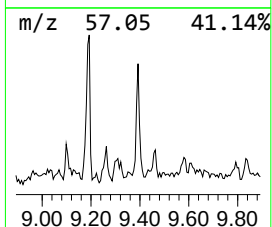
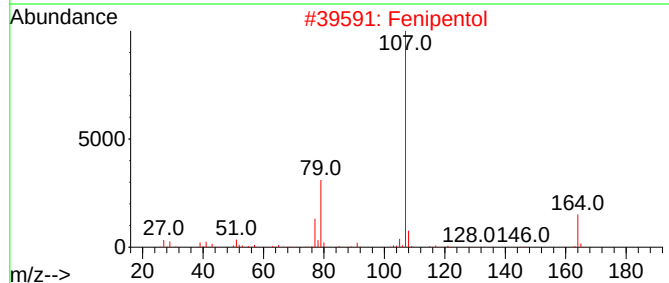
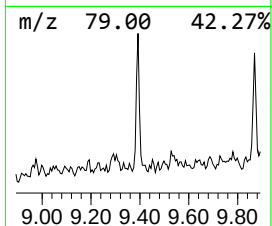
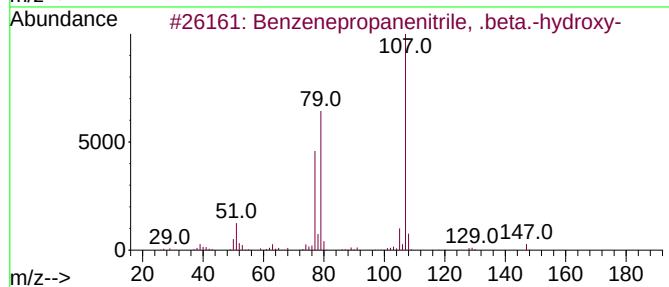
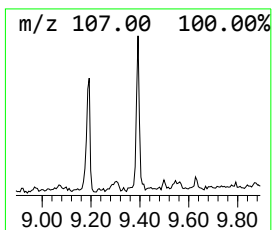
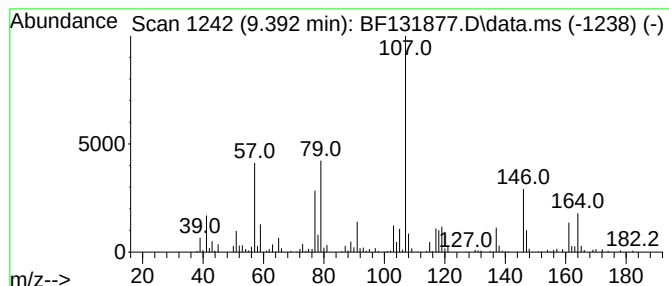
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 unknown9.392 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.392	3.80 ng	152448	Acenaphthene-d10	9.869

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzenepropanenitrile, .beta.-hy...	147	C9H9NO	017190-29-3	49
2			Fenipentol	164	C11H16O	000583-03-9	47
3			Benzenemethanol, .alpha.-propyl-	150	C10H14O	000614-14-2	47
4			Benzeneacetic acid, .alpha.-hydr...	166	C9H10O3	020698-91-3	43
5			2-Oxazolidinone, 4-methyl-5-phen...	177	C10H11NO2	028044-22-6	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF122922\
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Misc :
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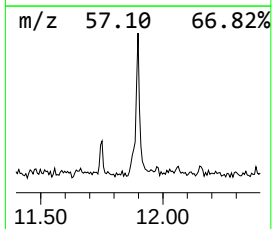
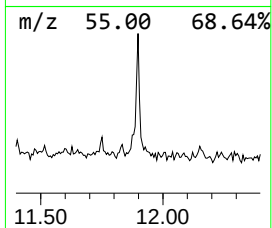
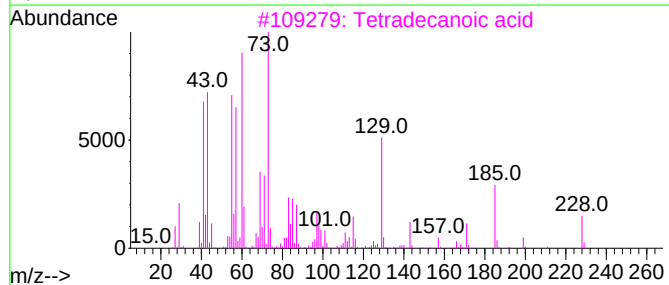
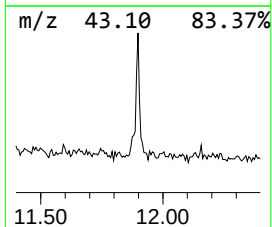
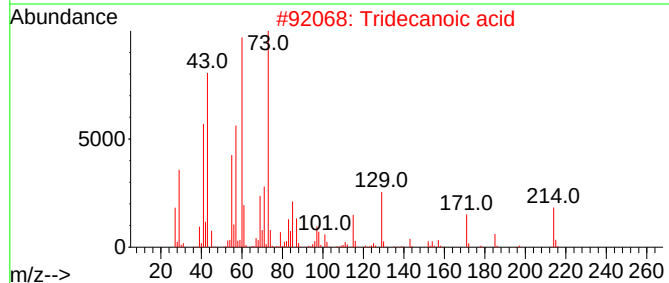
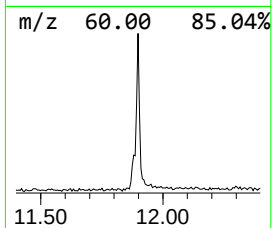
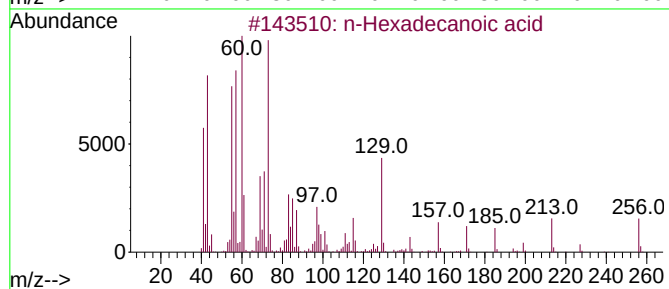
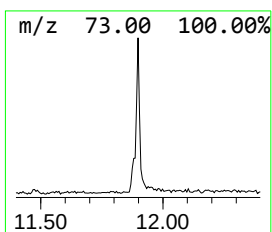
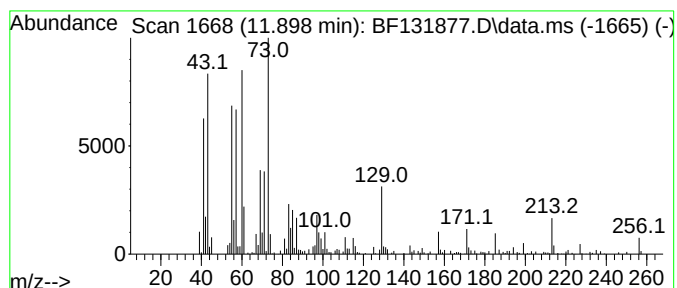
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 n-Hexadecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.898	5.08 ng	196855	Phenanthrene-d10	11.363	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2	Tridecanoic acid	214	C13H26O2	000638-53-9	87
3	Tetradecanoic acid	228	C14H28O2	000544-63-8	86
4	Pentadecanoic acid	242	C15H30O2	001002-84-2	86
5	Undecanoic acid	186	C11H22O2	000112-37-8	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF122922\
Data File : BF131877.D
Acq On : 29 Dec 2022 14:44
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Sample : N6232-08
Misc :
ALS Vial : 5 Sample Multiplier: 1

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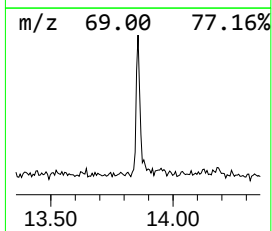
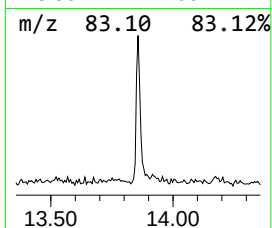
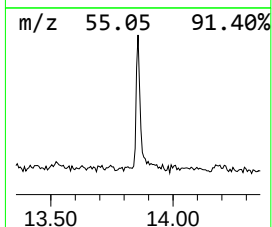
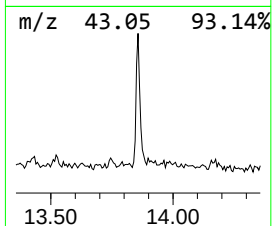
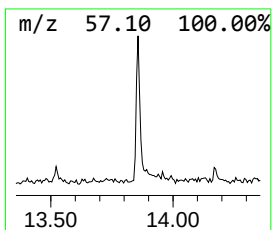
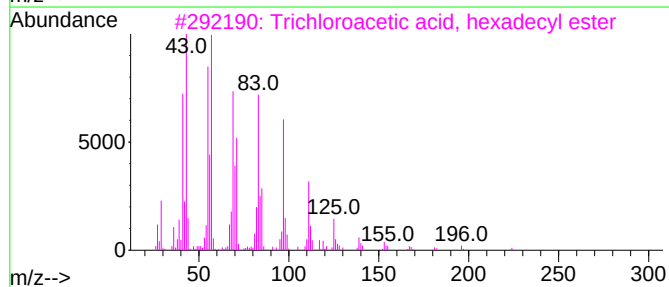
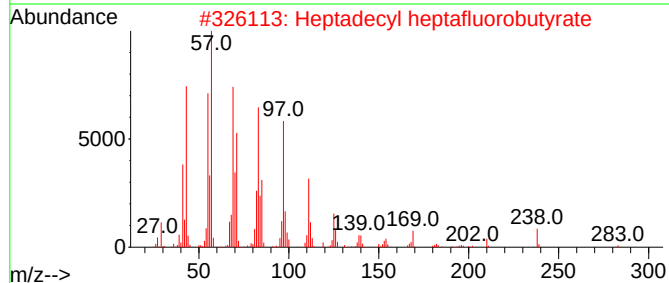
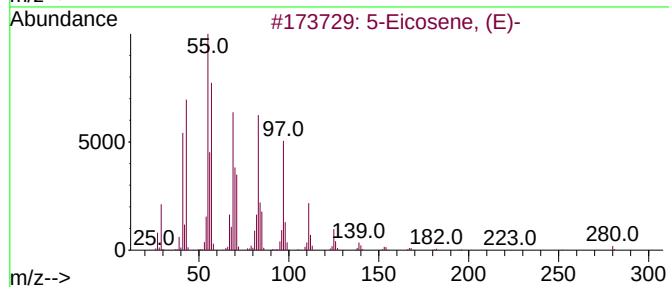
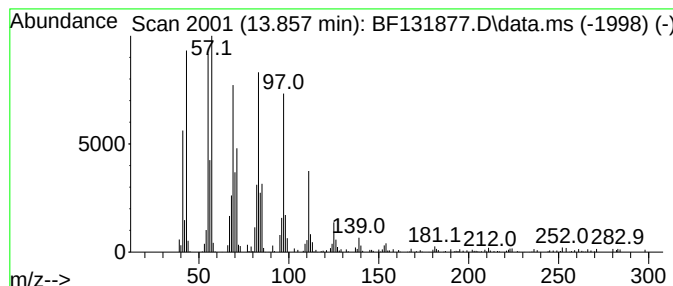
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 5-Eicosene, (E)- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.857	8.59 ng	218274	Chrysene-d12	14.016

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			5-Eicosene, (E)-	280	C20H40	074685-30-6	98
2			Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	94
3			Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	94
4			5-Octadecene, (E)-	252	C18H36	007206-21-5	93
5			1-Heneicosyl formate	340	C22H44O2	077899-03-7	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF122922\
Data File : BF131877.D
Acq On : 29 Dec 2022 14:44
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ALS Vial : 5 Sample Multiplier: 1

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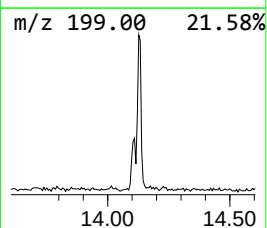
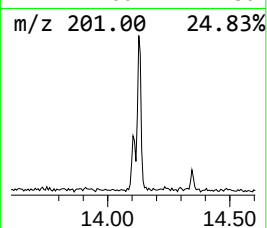
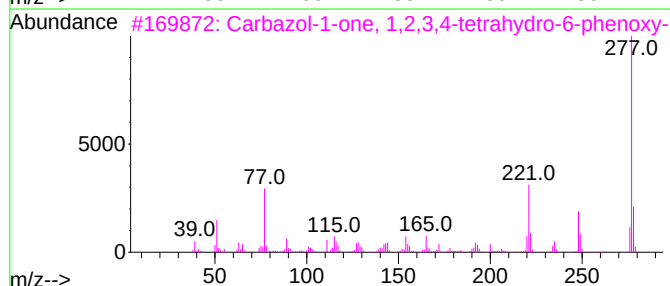
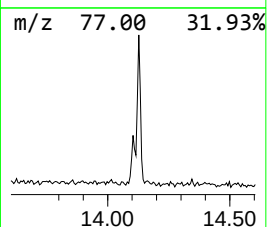
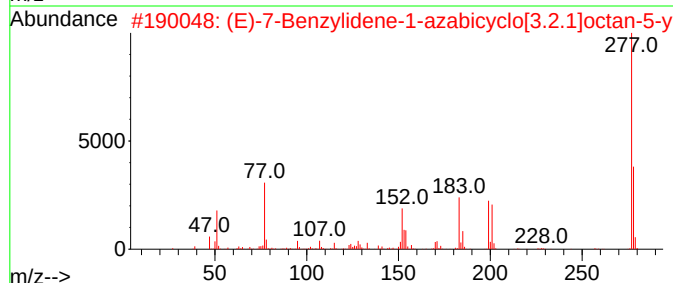
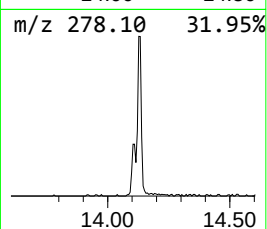
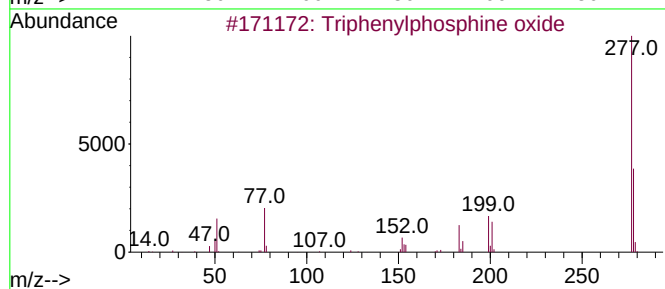
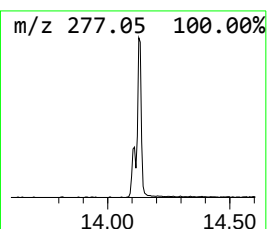
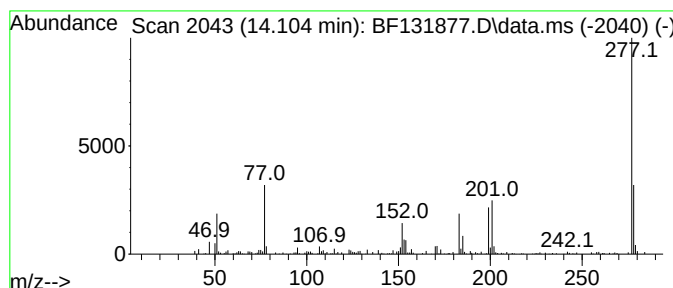
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Triphenylphosphine oxide Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.104	3.61 ng	91826	Chrysene-d12	14.016

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Triphenylphosphine oxide	278	C18H15OP	000791-28-6	93
2			(E)-7-Benzylidene-1-azabicyclo[3...]	293	C15H19NO3S	1000483-83-4	91
3			Carbazol-1-one, 1,2,3,4-tetrahyd...	277	C18H15NO2	299962-12-2	50
4			4-Nitro-4'-methyldiphenylsulfone	277	C13H11NO4S	004094-37-5	46
5			Formaldehyde, (triphenylphosphor...	304	C19H17N2P	015990-54-2	45



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF122922\
Data File : BF131877.D
Acq On : 29 Dec 2022 14:44
Operator : CG\JU
Sample : N6232-08
Misc :
ALS Vial : 5 Sample Multiplier: 1

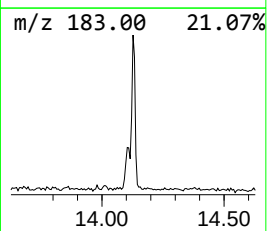
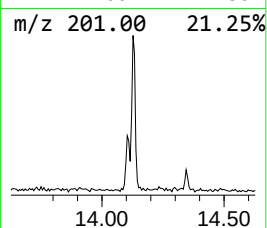
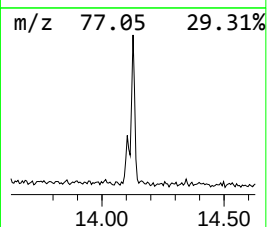
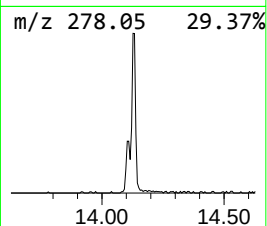
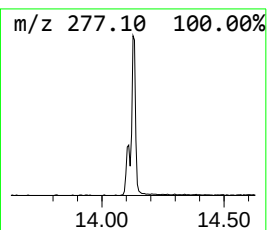
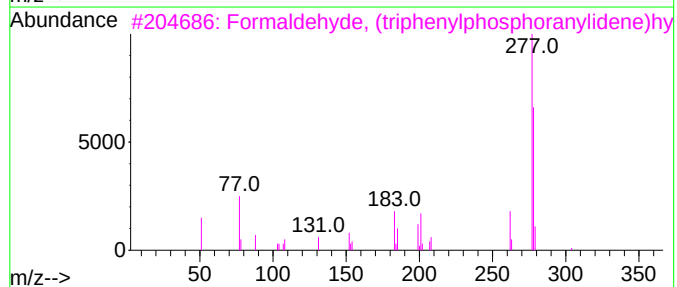
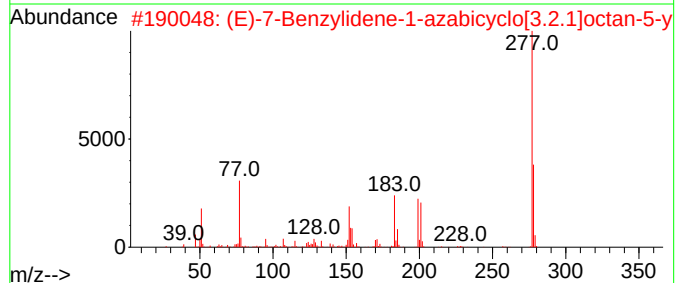
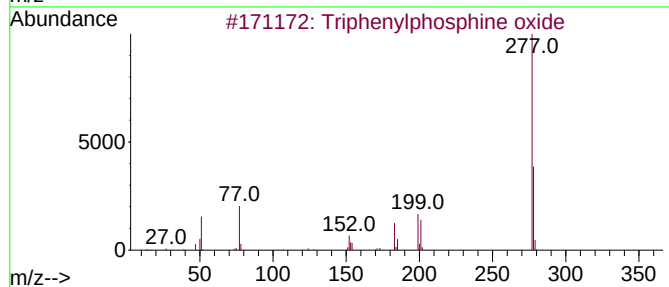
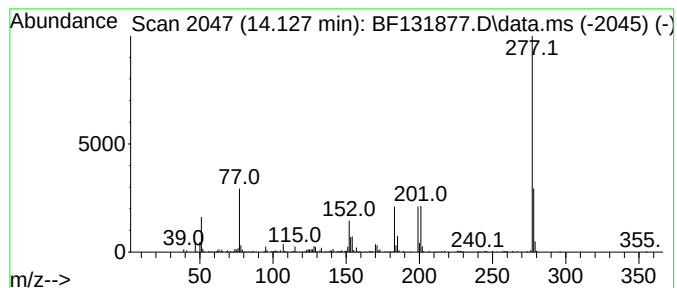
Instrument :
BNA_F
ClientSampleId :
MW-2

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF122422.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 (E)-7-Benzylidene-1-azabicy... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD		R.T.	
14.127	9.50 ng	241453	Chrysene-d12		14.016	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Triphenylphosphine oxide		278	C18H15OP	000791-28-6	98
2	(E)-7-Benzylidene-1-azabicyclo[3...		293	C15H19NO3S	1000483-83-4	91
3	Formaldehyde, (triphenylphosphor...		304	C19H17N2P	015990-54-2	50
4	Carbazol-1-one, 1,2,3,4-tetrahyd...		277	C18H15NO2	299962-12-2	47
5	2-Propenoic acid, 2-cyano-3-(3-p...		277	C18H15NO2	1000080-00-3	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF122922\
Data File : BF131877.D
Acq On : 29 Dec 2022 14:44
Operator : CG\JU
Sample : N6232-08
Misc :
ALS Vial : 5 Sample Multiplier: 1

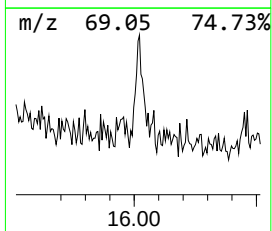
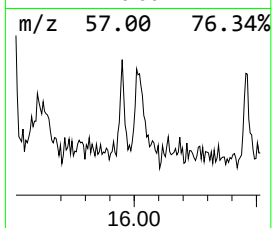
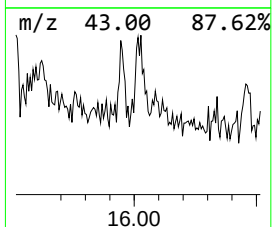
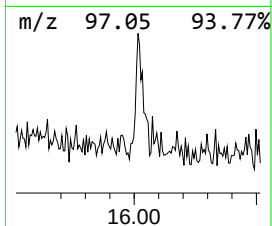
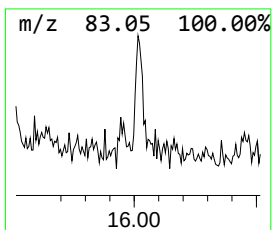
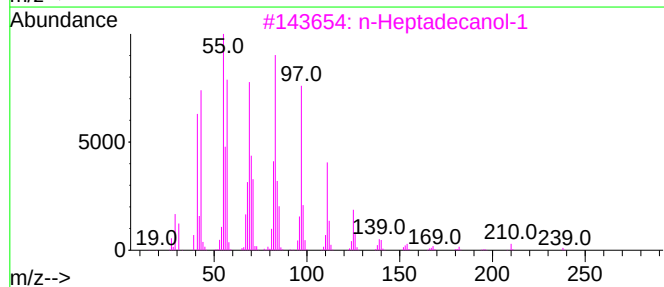
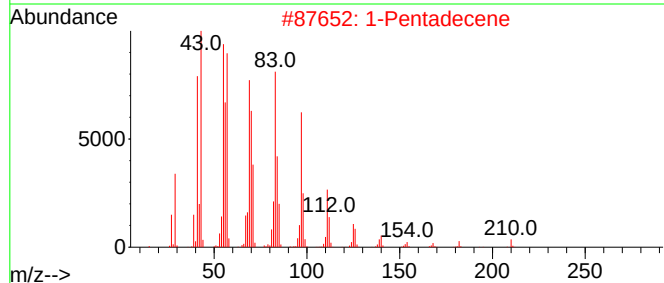
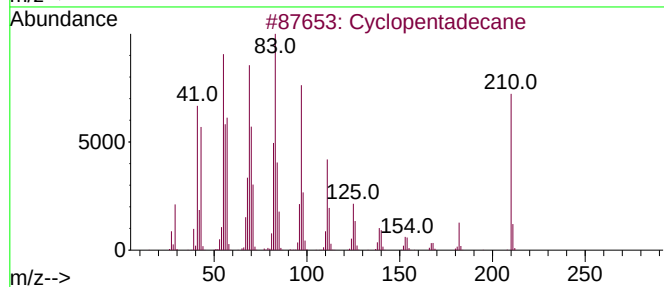
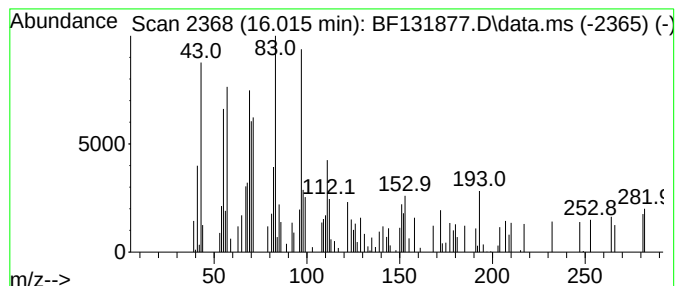
Instrument :
BNA_F
ClientSampleId :
MW-2

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF122422.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Cyclopentadecane Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.	
16.015	2.05 ng	47906	Perylene-d12	15.498	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentadecane	210	C15H30	000295-48-7	89
2	1-Pentadecene	210	C15H30	013360-61-7	89
3	n-Heptadecanol-1	256	C17H36O	001454-85-9	68
4	Z-5-Nonadecene	266	C19H38	1000131-11-8	68
5	n-Pentadecanol	228	C15H32O	000629-76-5	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF122922\
Data File : BF131877.D
Acq On : 29 Dec 2022 14:44
Operator : CG\JU
Sample : N6232-08
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MW-2

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF122422.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
2-Pentanone, 4-...	5.016	10.2	ng	277831	1	6.816	545297	20.0
Hexanoic acid, ...	7.698	30.6	ng	1114150	2	8.104	728943	20.0
(+)-2-Bornanone	7.851	2.4	ng	88294	2	8.104	728943	20.0
1-methyl-1-indanol	8.328	3.4	ng	123904	2	8.104	728943	20.0
Phenol, p-tert-...	8.681	2.6	ng	93815	2	8.104	728943	20.0
unknown8.863	8.863	2.4	ng	85564	2	8.104	728943	20.0
unknown9.304	9.304	2.8	ng	112677	3	9.869	803359	20.0
unknown9.392	9.392	3.8	ng	152448	3	9.869	803359	20.0
n-Hexadecanoic ...	11.898	5.1	ng	196855	4	11.363	775469	20.0
5-Eicosene, (E)-	13.857	8.6	ng	218274	5	14.016	508107	20.0
Triphenylphosph...	14.104	3.6	ng	91826	5	14.016	508107	20.0
(E)-7-Benzylide...	14.127	9.5	ng	241453	5	14.016	508107	20.0
Cyclopentadecane	16.015	2.0	ng	47906	6	15.498	467784	20.0