

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF043020\
 Data File : BF120245.D
 Acq On : 30 Apr 2020 17:55
 Operator : MA/SJ
 Sample : L2451-01
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MW-1

Integration Parameters: rteint.p
 Integrator: RTE
 Smoothing : ON
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0
 Filtering: 5
 Min Area: 3 % of largest Peak
 Max Peaks: 100
 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-BF040820.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.251	535	538	553	rBV	1456886	1486427	36.12%	6.116%
2	6.298	713	716	728	rBV	1064165	963984	23.43%	3.966%
3	6.645	772	775	781	rBV	682954	556360	13.52%	2.289%
4	6.804	798	802	805	rBV	2682784	2213295	53.79%	9.107%
5	7.216	865	872	875	rBV	2172545	2032033	49.38%	8.361%
6	7.934	990	994	997	rBV	958394	707125	17.18%	2.910%
7	8.175	1031	1035	1043	rBV	134441	147309	3.58%	0.606%
8	8.663	1113	1118	1128	rBV	299901	336993	8.19%	1.387%
9	8.733	1128	1130	1141	rVB	40099	50575	1.23%	0.208%
10	8.845	1146	1149	1153	rVV	105613	89009	2.16%	0.366%
11	8.898	1155	1158	1162	rVB	59818	47057	1.14%	0.194%
12	9.016	1174	1178	1181	rBV	4265315	3465857	84.23%	14.261%
13	9.416	1242	1246	1249	rBV	301919	253002	6.15%	1.041%
14	9.692	1289	1293	1296	rBV	954403	839898	20.41%	3.456%
15	10.139	1366	1369	1375	rVB2	106006	99010	2.41%	0.407%
16	10.416	1410	1416	1423	rBV	68267	85899	2.09%	0.353%
17	10.486	1423	1428	1431	rBV	2362627	2361826	57.40%	9.718%
18	10.922	1498	1502	1515	rBV	505025	528418	12.84%	2.174%
19	11.174	1542	1545	1549	rBV	1076169	849338	20.64%	3.495%
20	11.722	1634	1638	1645	rBV2	246983	271579	6.60%	1.117%
21	12.504	1768	1771	1782	rBV	92217	110226	2.68%	0.454%
22	12.651	1793	1796	1809	rBV	205936	221381	5.38%	0.911%
23	12.774	1811	1817	1820	rBV	3936030	4114853	100.00%	16.931%
24	13.674	1966	1970	1983	rBV	153528	227072	5.52%	0.934%
25	13.815	1990	1994	1998	rBV	967989	826651	20.09%	3.401%
26	14.298	2072	2076	2086	rBV2	37545	66610	1.62%	0.274%
27	14.904	2176	2179	2185	rBV2	42534	48145	1.17%	0.198%
28	15.221	2228	2233	2236	rBV	667383	797838	19.39%	3.283%
29	15.257	2236	2239	2244	rVB	43704	54294	1.32%	0.223%
30	15.651	2302	2306	2313	rBV2	29857	45524	1.11%	0.187%
31	16.956	2521	2528	2532	rBV6	38151	84452	2.05%	0.347%
32	17.062	2540	2546	2552	rVB5	43013	80939	1.97%	0.333%
33	18.456	2775	2783	2792	rBV6	64813	167609	4.07%	0.690%
34	18.609	2801	2809	2816	rBV6	26067	72644	1.77%	0.299%

LSC Area Percent Report

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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-BF040820.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

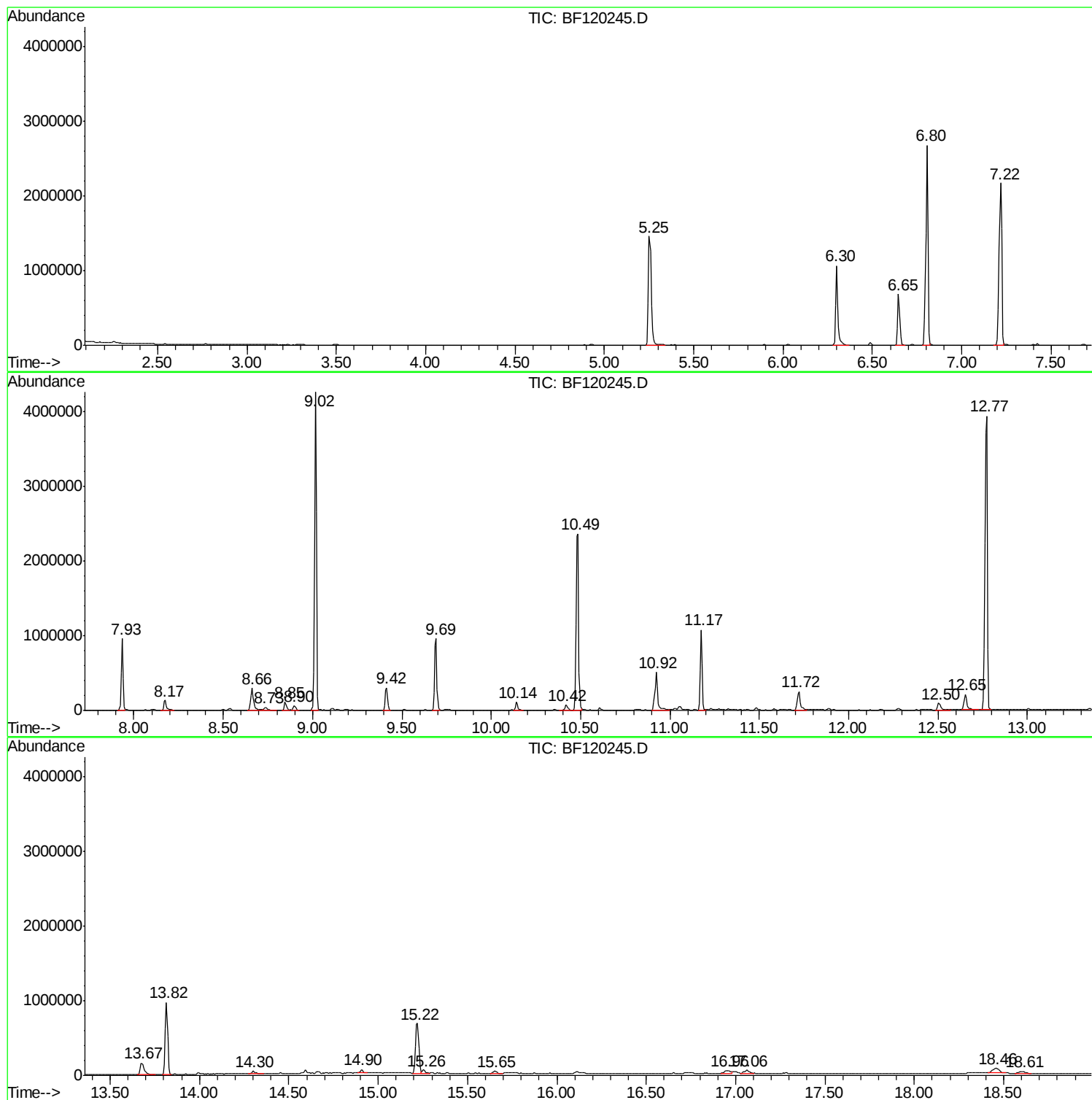
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF040820.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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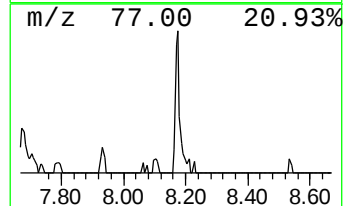
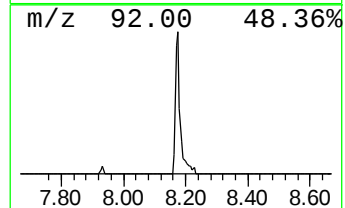
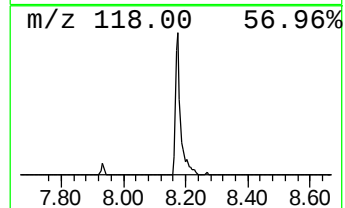
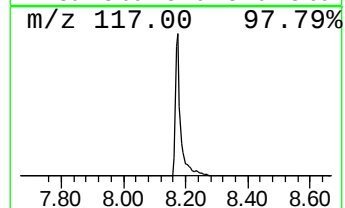
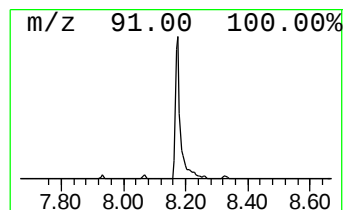
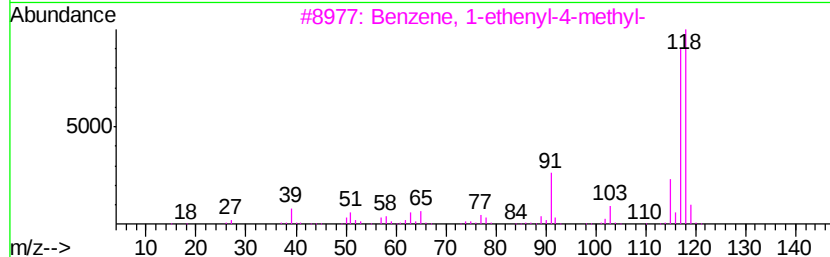
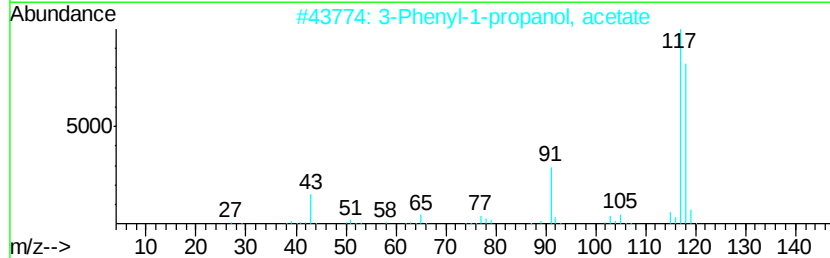
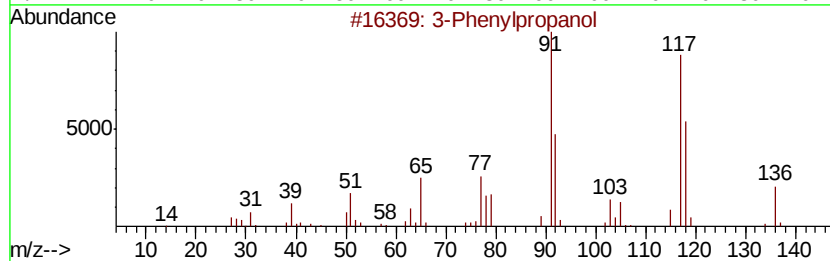
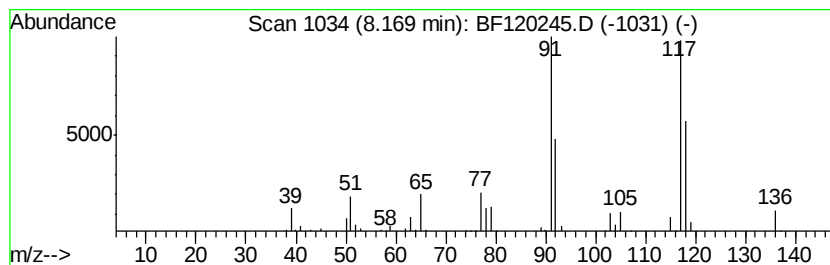
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 3-Phenylpropanol Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.17	4.17 ng	147309	Naphthalene-d8	7.93

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Phenylpropanol	136	C9H12O	000122-97-4	96
2		3-Phenyl-1-propanol, acetate	178	C11H14O2	000122-72-5	80
3		Benzene, 1-ethenyl-4-methyl-	118	C9H10	000622-97-9	64
4		Tetracyclo[3.3.1.0(2,8).0(4,6)]-	118	C9H10	1000191-13-7	53
5		1-Methoxy-3-phenylpropane	150	C10H14O	002046-33-5	50



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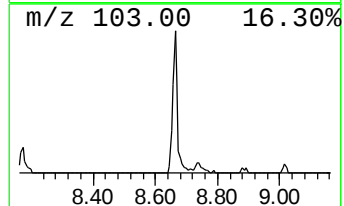
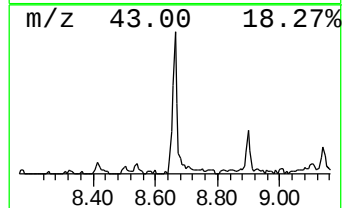
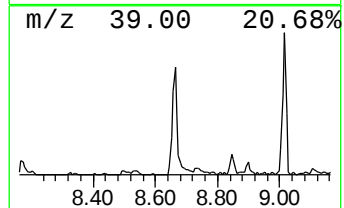
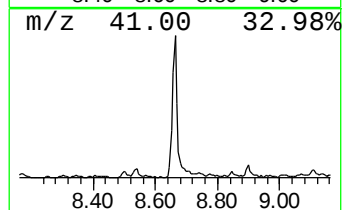
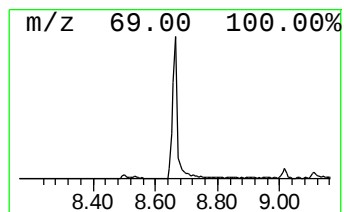
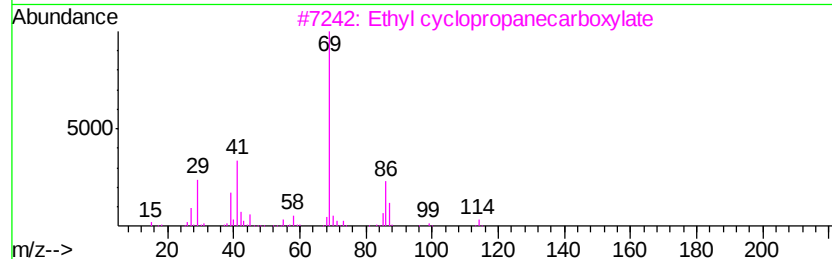
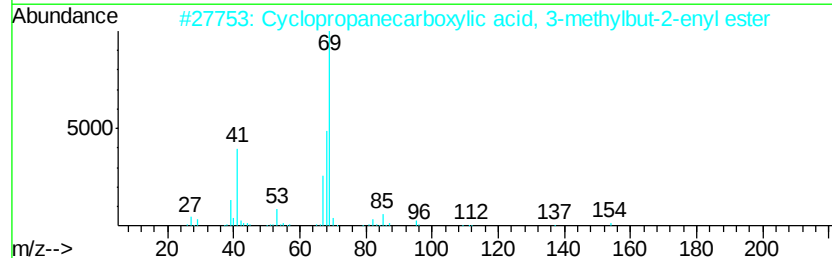
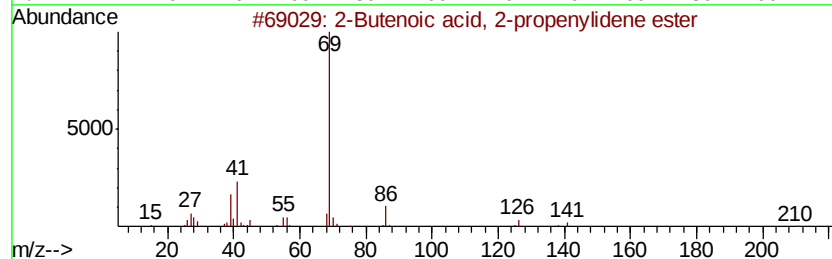
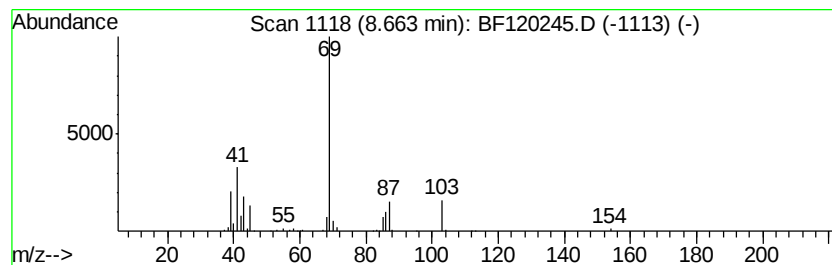
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 unknown8.66 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.66	9.53 ng	336993	Napthalene-d8	7.93

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Butenoic acid, 2-propenylidene...	210	C11H14O4	055030-70-1	45
2		Cyclopropanecarboxylic acid, 3-m...	154	C9H14O2	1000299-37-4	40
3		Ethyl cyclopropanecarboxylate	114	C6H10O2	004606-07-9	38
4		.alpha.-Hydroxyisocaproic acid	132	C6H12O3	000498-36-2	9
5		(+)-3-hydroxybutyric acid, trif...	200	C6H7F3O4	1000374-31-1	9



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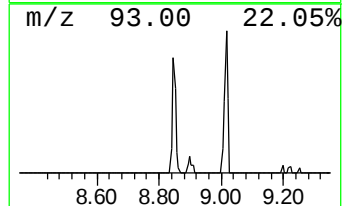
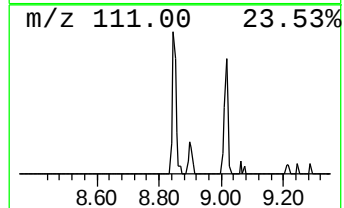
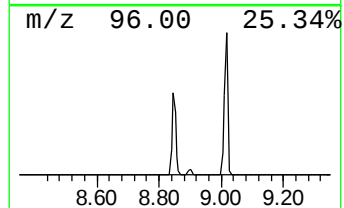
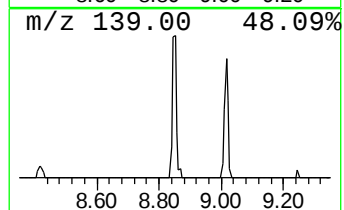
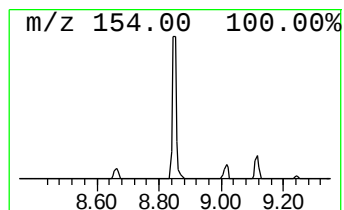
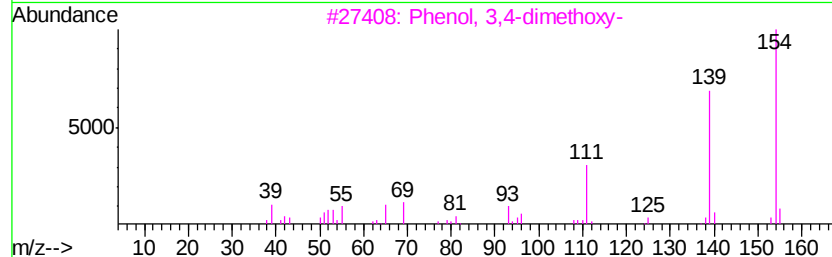
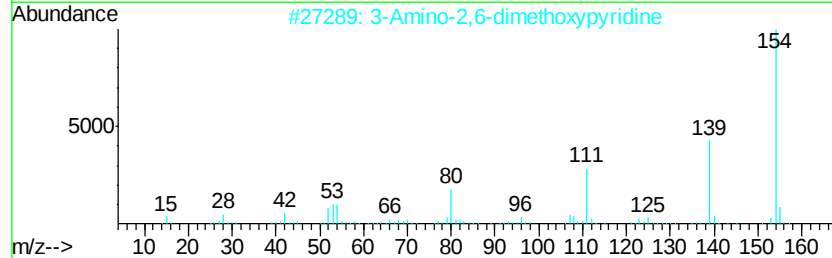
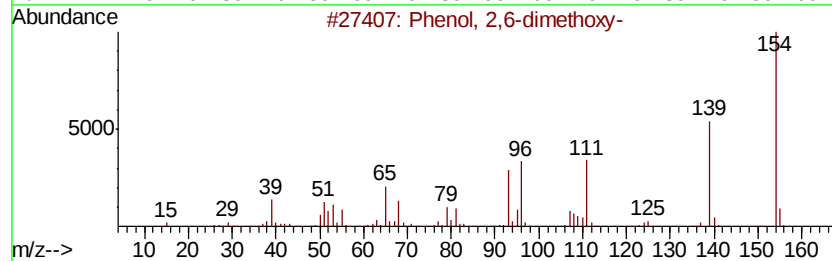
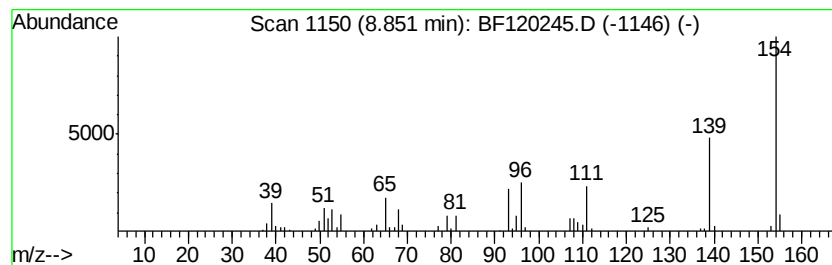
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Phenol, 2,6-dimethoxy- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.		
8.85	2.12 ng	89009	Acenaphthene-d10	9.69		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 2,6-dimethoxy-	154	C8H10O3	000091-10-1	97	
2	3-Amino-2,6-dimethoxypyridine	154	C7H10N2O2	028020-37-3	70	
3	Phenol, 3,4-dimethoxy-	154	C8H10O3	002033-89-8	59	
4	Octahydro-1H-pyrido(1,2-c)pyrimi...	154	C8H14N2O	015932-63-5	47	
5	Phenol, 3,5-dimethoxy-	154	C8H10O3	000500-99-2	32	



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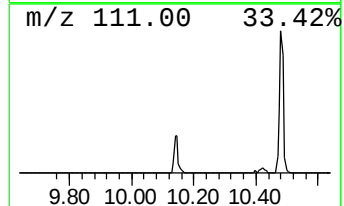
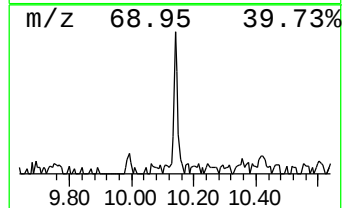
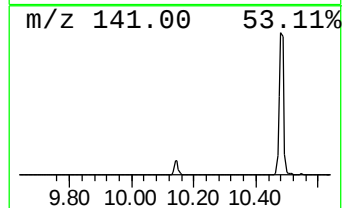
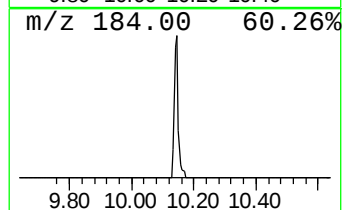
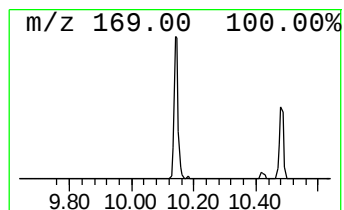
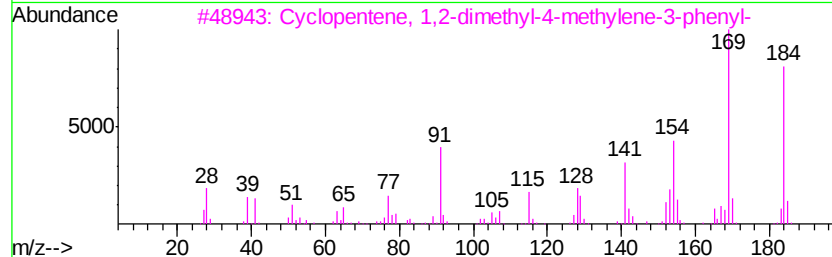
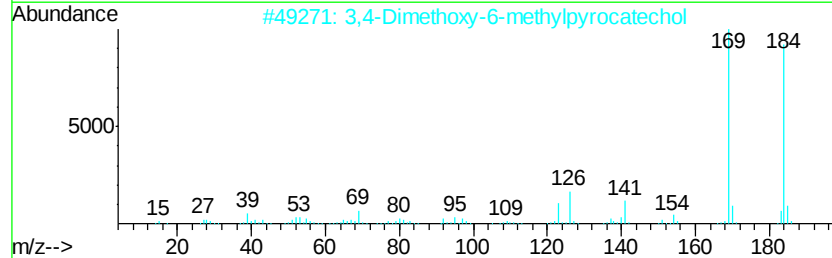
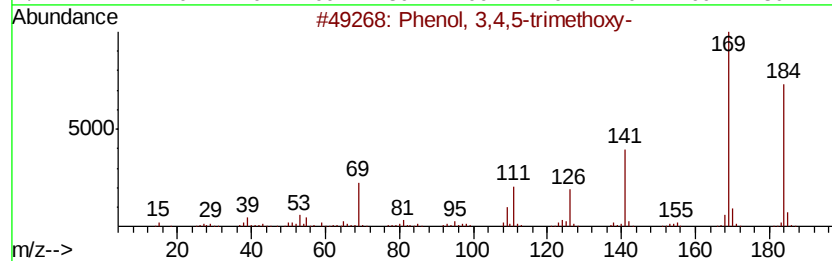
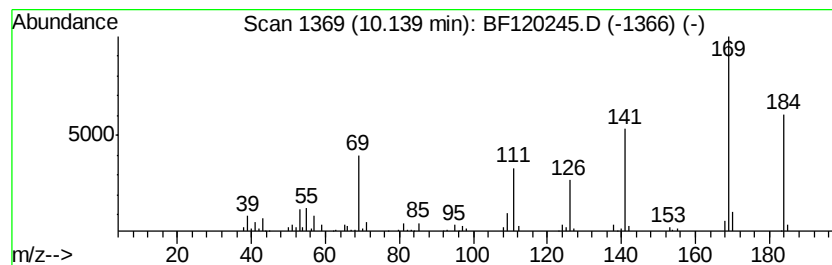
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Phenol, 3,4,5-trimethoxy- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.14	2.36 ng	99010	Acenaphthene-d10	9.69
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Phenol, 3,4,5-trimethoxy-	184 C9H12O4	000642-71-7	96
2	3,4-Dimethoxy-6-methylpyrocatechol	184 C9H12O4	054826-79-8	50
3	Cyclopentene, 1,2-dimethyl-4-met...	184 C14H16	1000152-22-4	50
4	3-Thiopheneacetic acid, 5-acetyl-	184 C8H8O3S	1000337-56-9	49
5	2(1H)-Pyridinethione, 1-ethyl-3-...	169 C8H11NOS	024207-15-6	47



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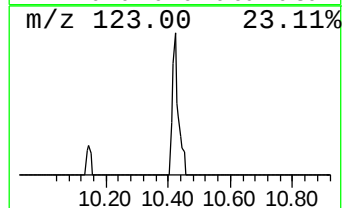
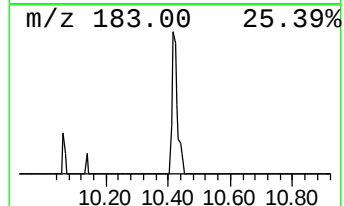
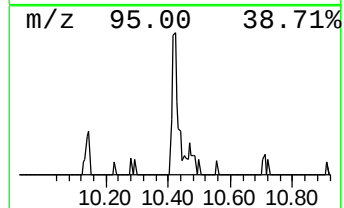
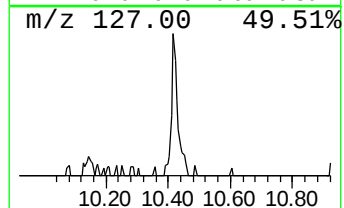
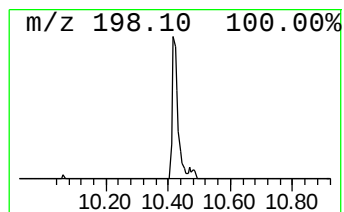
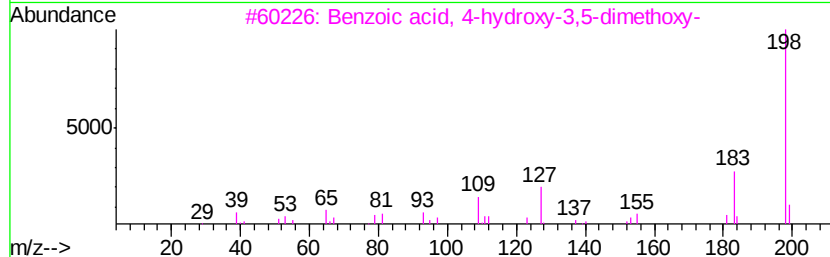
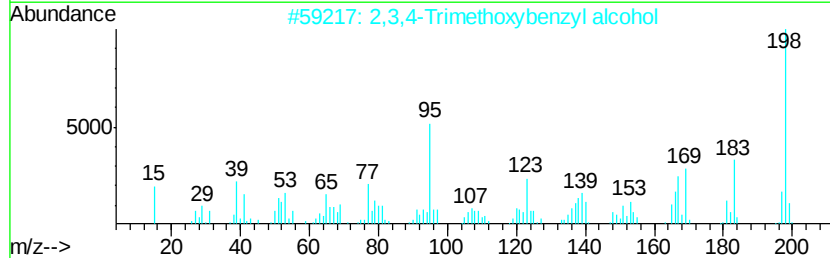
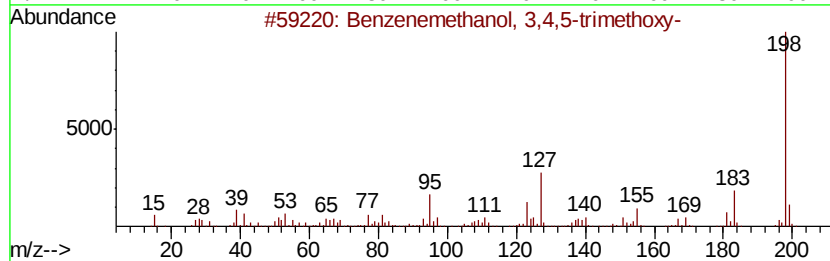
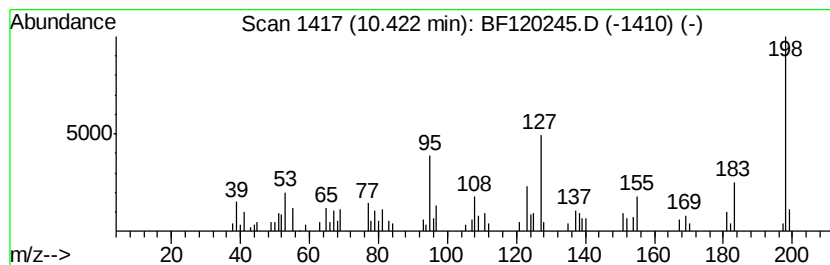
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Benzenemethanol, 3,4,5-trim... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.42	2.05 ng	85899	Acenaphthene-d10	9.69

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzenemethanol, 3,4,5-trimethoxy-	198	C10H14O4	003840-31-1	95
2		2,3,4-Trimethoxybenzyl alcohol	198	C10H14O4	071989-96-3	76
3		Benzoic acid, 4-hydroxy-3,5-dime...	198	C9H10O5	000530-57-4	49
4		3-Hydroxy-4,5-dimethoxybenzoic acid	198	C9H10O5	001916-08-1	47
5		Benzene, 1,1'-oxybis[4-methyl-	198	C14H14O	001579-40-4	46



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF043020\
Data File : BF120245.D
Acq On : 30 Apr 2020 17:55
Operator : MA/SJ
Sample : L2451-01
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MW-1

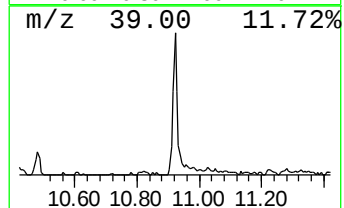
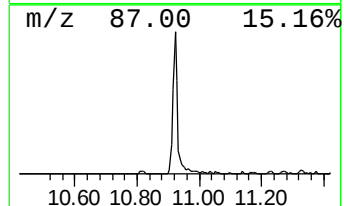
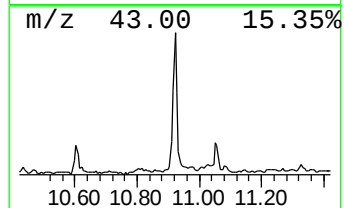
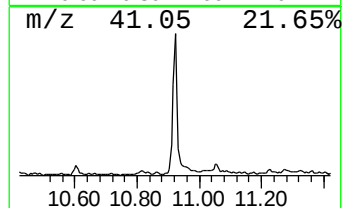
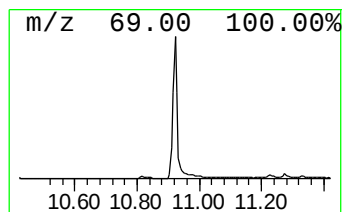
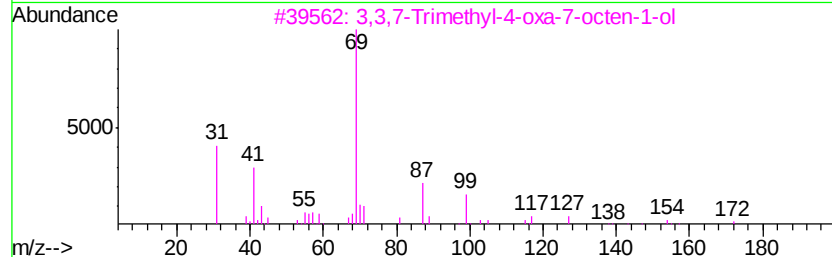
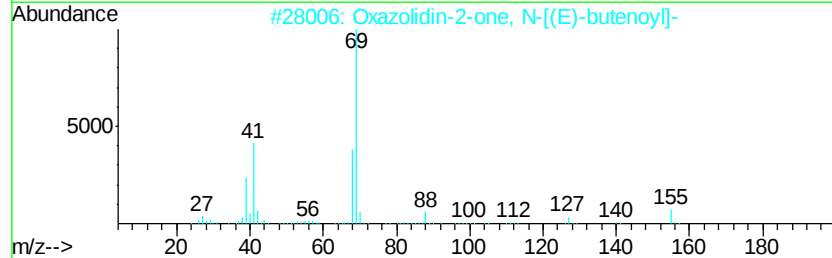
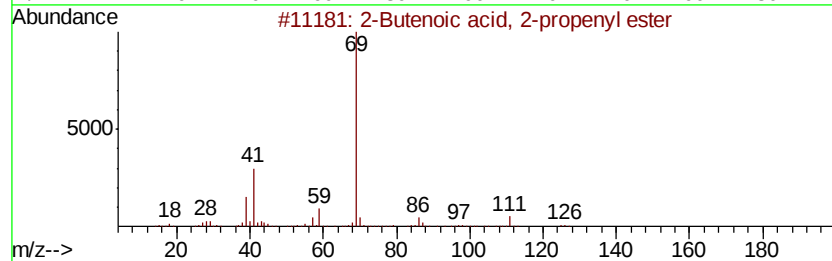
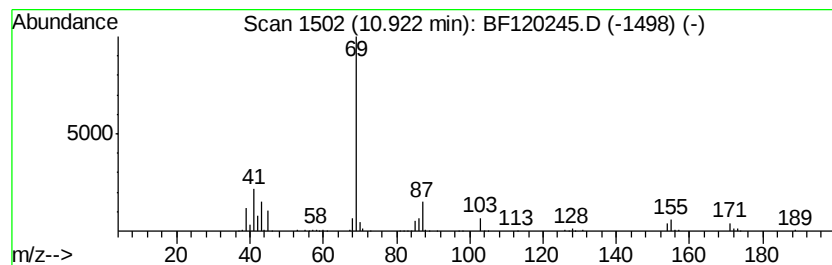
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 unknown10.92 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.92	12.44 ng	528418	Phenanthrene-d10	11.17

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Butenoic acid, 2-propenyl ester	126	C7H10O2	020474-93-5	40
2		Oxazolidin-2-one, N-[(E)-butenoyl]-	155	C7H9NO3	1000156-45-5	40
3		3,3,7-Trimethyl-4-oxa-7-octen-1-ol	172	C10H20O2	069161-47-3	38
4		Cyclopropanecarboxylic acid, isob...	142	C8H14O2	064969-79-5	9
5		Ethyl cyclopropanecarboxylate	114	C6H10O2	004606-07-9	9



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF043020\
Data File : BF120245.D
Acq On : 30 Apr 2020 17:55
Operator : MA/SJ
Sample : L2451-01
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MW-1

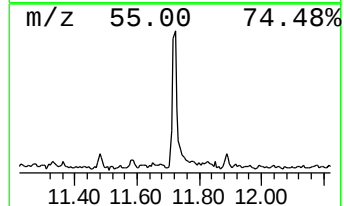
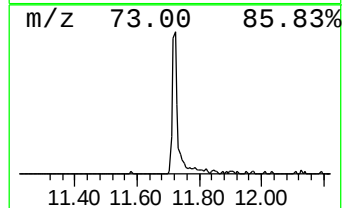
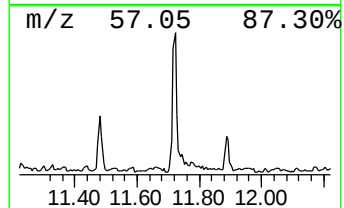
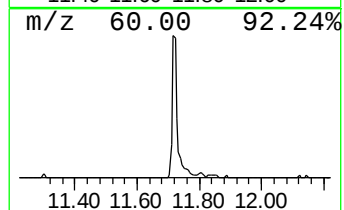
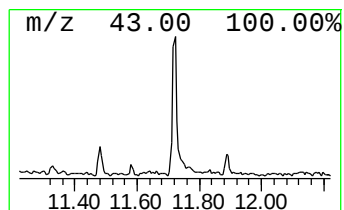
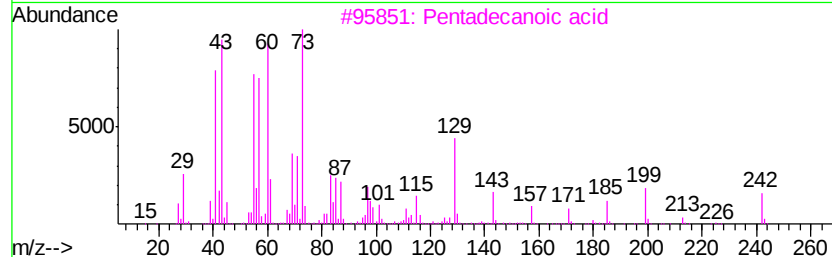
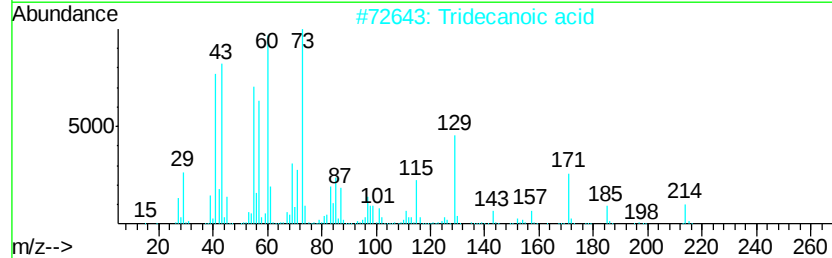
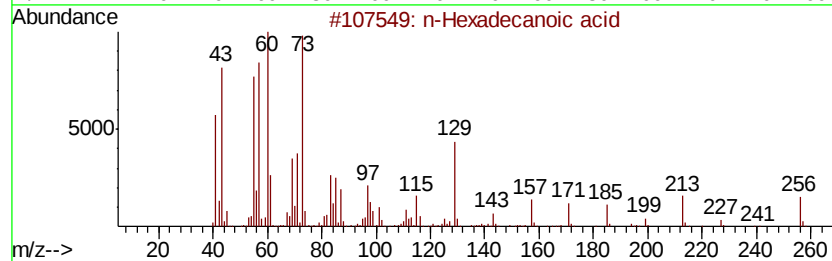
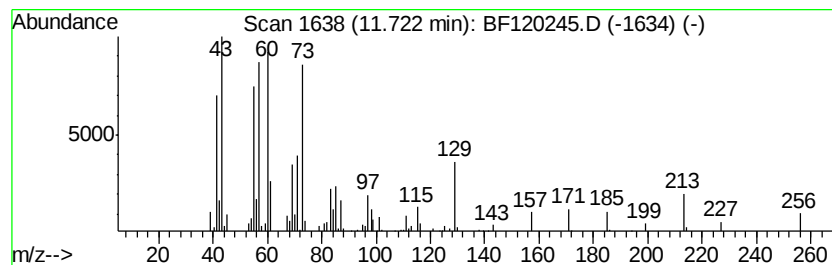
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 n-Hexadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.72	6.40 ng	271579	Phenanthrene-d10	11.17

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	90
3	Pentadecanoic acid	242	C15H30O2	001002-84-2	87
4	n-Decanoic acid	172	C10H20O2	000334-48-5	64
5	Oxalic acid, dodecyl propyl ester	300	C17H32O4	1000309-26-5	25



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF043020\
Data File : BF120245.D
Acq On : 30 Apr 2020 17:55
Operator : MA/SJ
Sample : L2451-01
Misc :
ALS Vial : 14 Sample Multiplier: 1

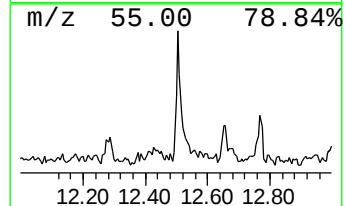
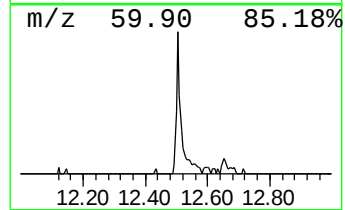
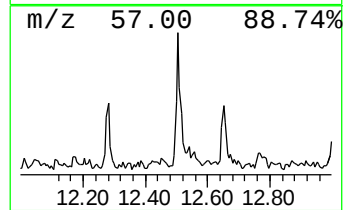
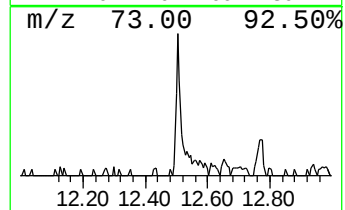
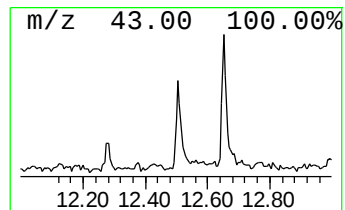
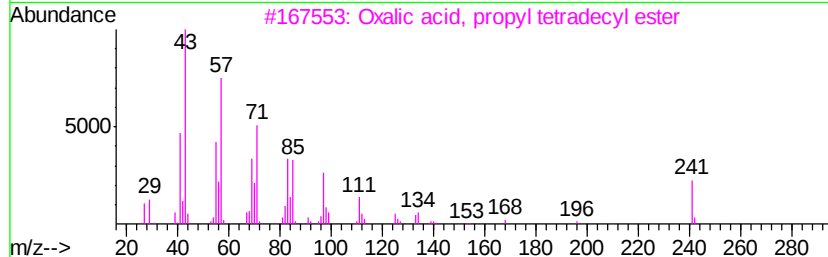
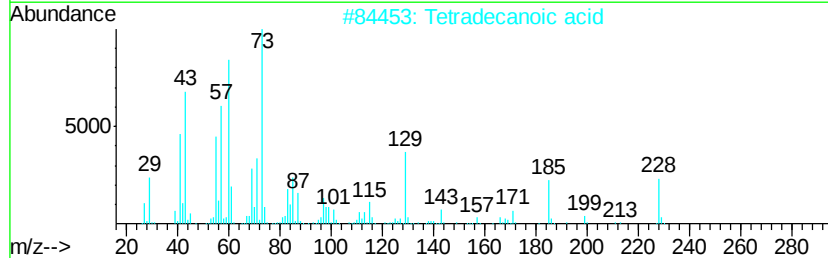
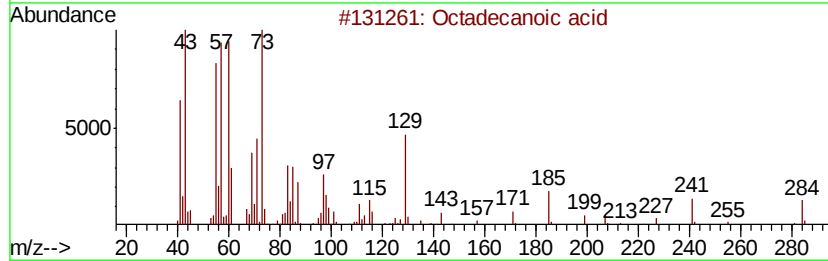
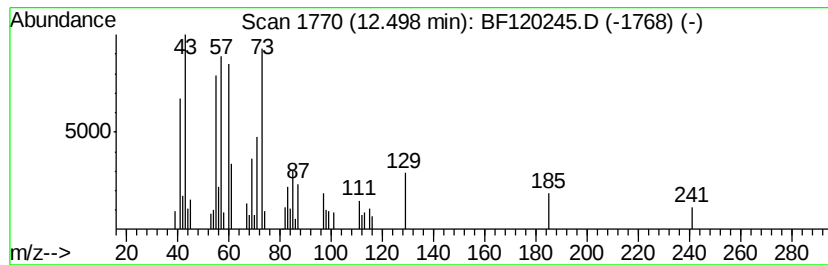
Instrument :
BNA_F
ClientSampleId :
MW-1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF040820.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Octadecanoic acid Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD		R.T.
12.50	2.67 ng	110226	Chrysene-d12		13.82
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecanoic acid	284	C18H36O2	000057-11-4	93
2	Tetradecanoic acid	228	C14H28O2	000544-63-8	64
3	Oxalic acid, propyl tetradecyl e...	328	C19H36O4	1000309-26-7	38
4	Oxalic acid, allyl octadecyl ester	382	C23H42O4	1000309-24-5	20
5	Oxalic acid, cyclobutyl hexadecy...	368	C22H40O4	1000309-70-6	18



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF043020\
Data File : BF120245.D
Acq On : 30 Apr 2020 17:55
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Sample : L2451-01
Misc :
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Instrument :
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ClientSampleId :
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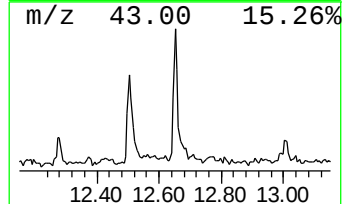
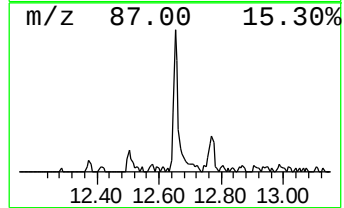
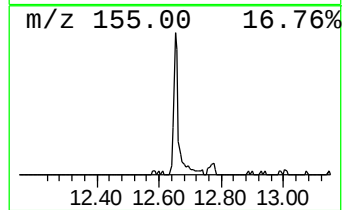
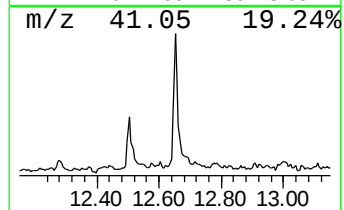
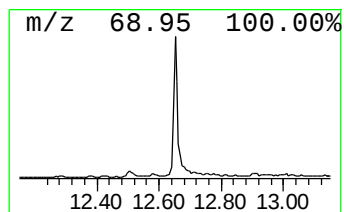
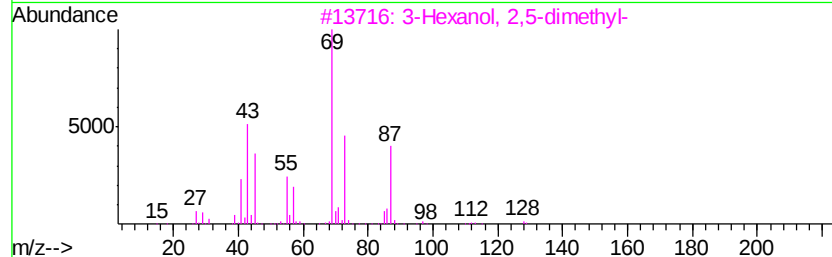
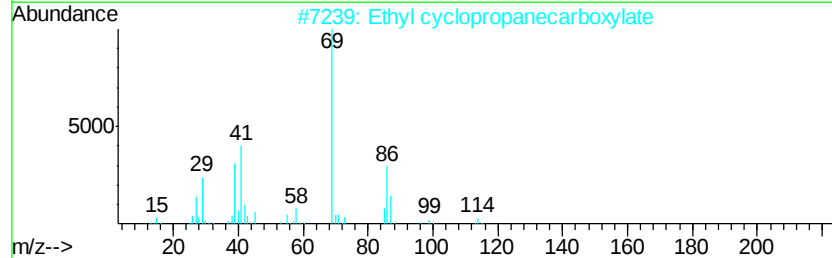
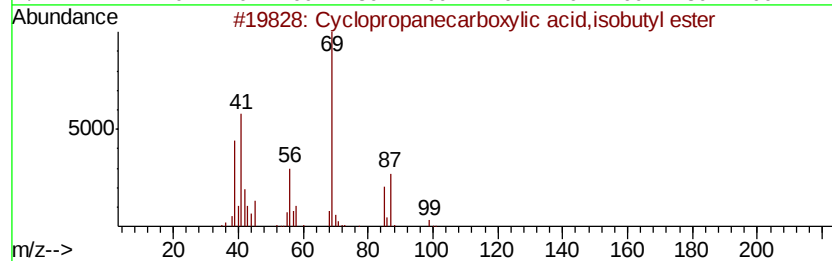
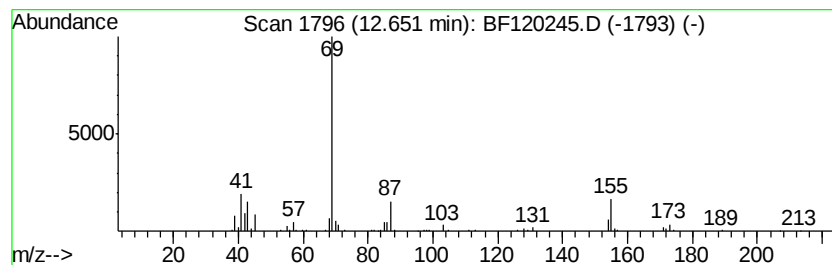
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Cyclopropanecarboxylic acid... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.65	5.36 ng	221381	Chrysene-d12	13.82

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopropanecarboxylic acid, isob...	142	C8H14O2	064969-79-5	59
2		Ethyl cyclopropanecarboxylate	114	C6H10O2	004606-07-9	38
3		3-Hexanol, 2,5-dimethyl-	130	C8H18O	019550-07-3	38
4		2-Butenoic acid, hexyl ester	170	C10H18O2	019089-92-0	38
5		2-Propenoic acid, 2-methyl-, 1-m...	128	C7H12O2	004655-34-9	36



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF043020\
Data File : BF120245.D
Acq On : 30 Apr 2020 17:55
Operator : MA/SJ
Sample : L2451-01
Misc :
ALS Vial : 14 Sample Multiplier: 1

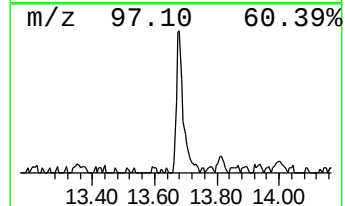
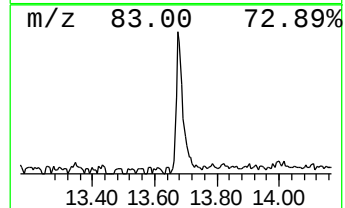
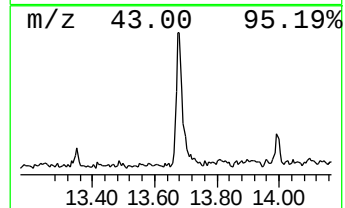
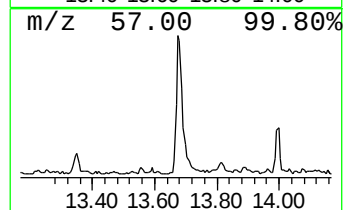
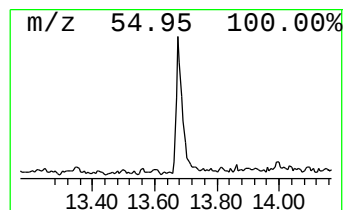
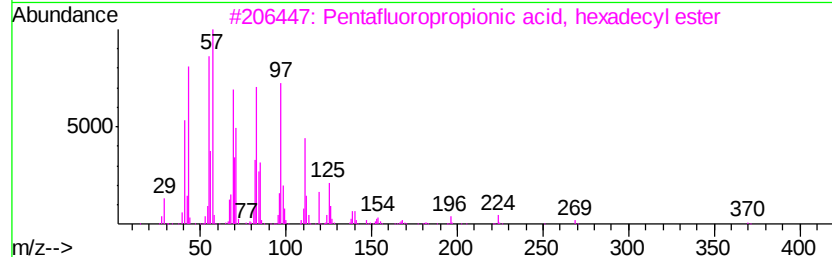
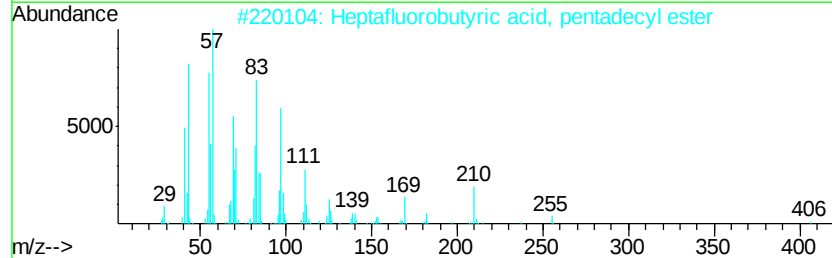
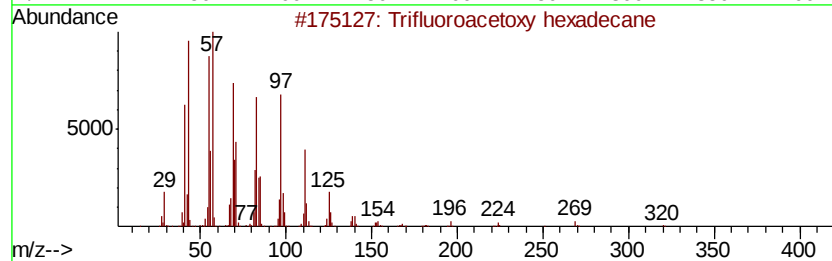
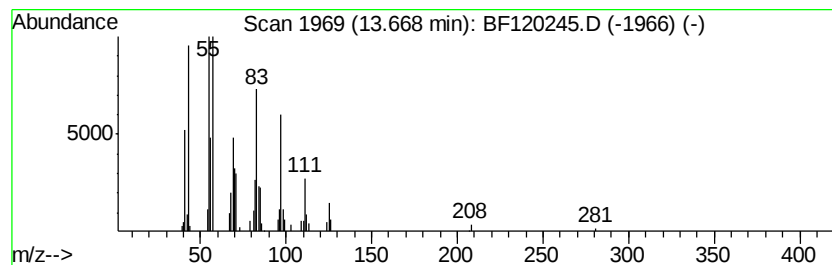
Instrument :
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF040820.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Trifluoroacetoxy hexadecane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.		
13.67	5.49 ng	227072	Chrysene-d12	13.82		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Trifluoroacetoxy hexadecane	338	C18H33F3O2	006222-03-3	91
2		Heptafluorobutyric acid, pentadec...	424	C19H31F7O2	959261-23-5	91
3		Pentafluoropropionic acid, hexad...	388	C19H33F5O2	006222-07-7	91
4		Heptafluorobutyric acid, hexadec...	438	C20H33F7O2	006385-15-5	91
5		2- Chloropropionic acid, octadec...	360	C21H41ClO2	088104-31-8	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF043020\
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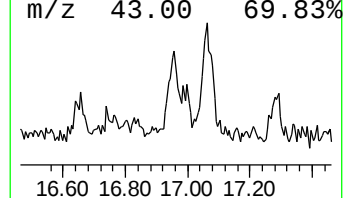
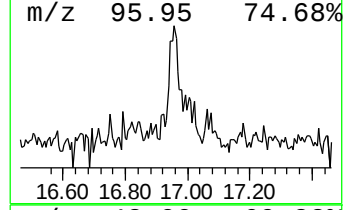
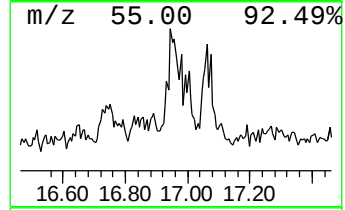
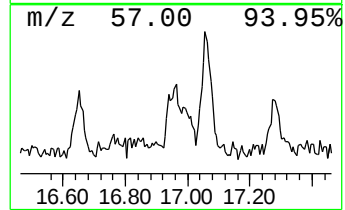
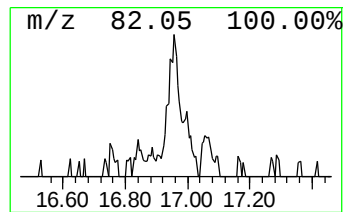
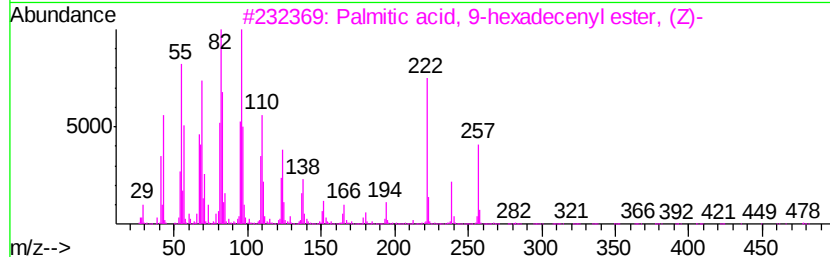
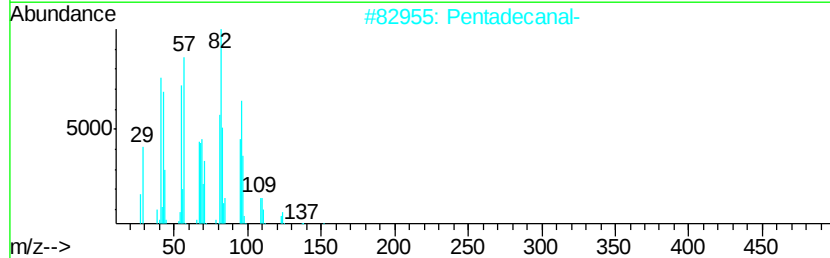
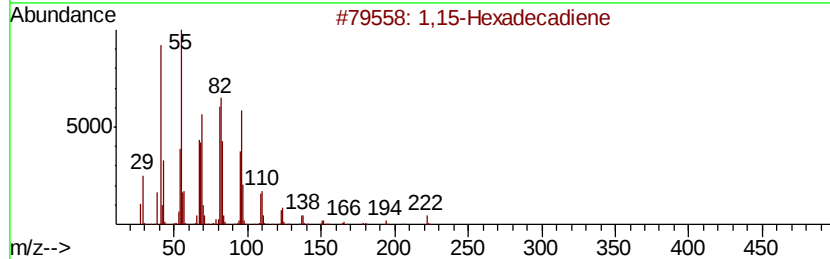
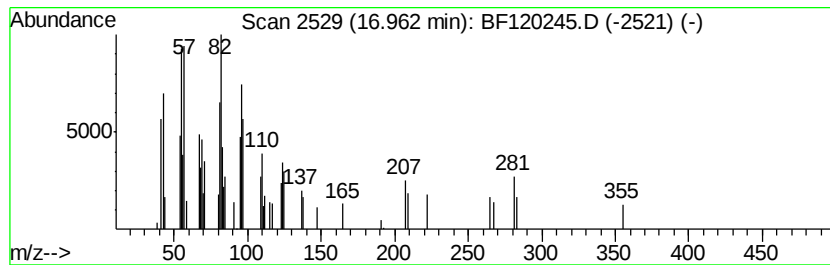
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 1,15-Hexadecadiene Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.96	2.12 ng	84452	Perylene-d12	15.22
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	1,15-Hexadecadiene	222 C16H30	021964-51-2	64
2	Pentadecanal-	226 C15H30O	002765-11-9	52
3	Palmitic acid, 9-hexadecenyl est...	478 C32H62O2	022393-90-4	50
4	9-Octadecen-1-ol, (E)-	268 C18H36O	000506-42-3	49
5	Bicyclo[4.1.0]heptane, 7-butyl-	152 C11H20	018645-10-8	46



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF043020\
Data File : BF120245.D
Acq On : 30 Apr 2020 17:55
Operator : MA/SJ
Sample : L2451-01
Misc :
ALS Vial : 14 Sample Multiplier: 1

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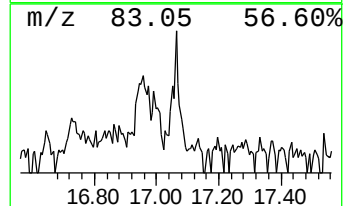
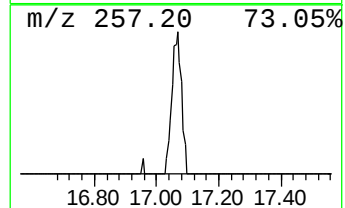
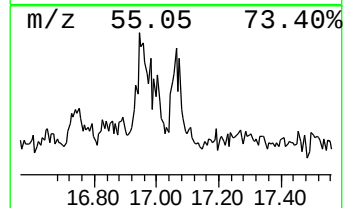
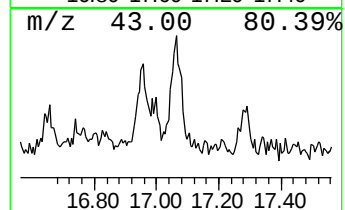
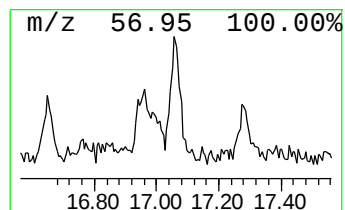
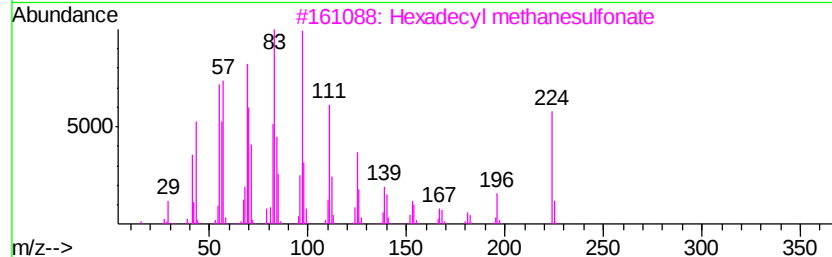
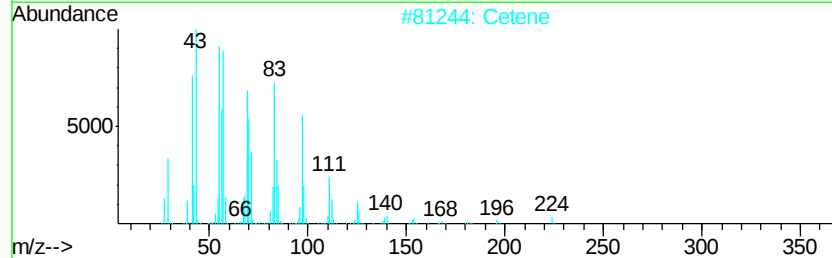
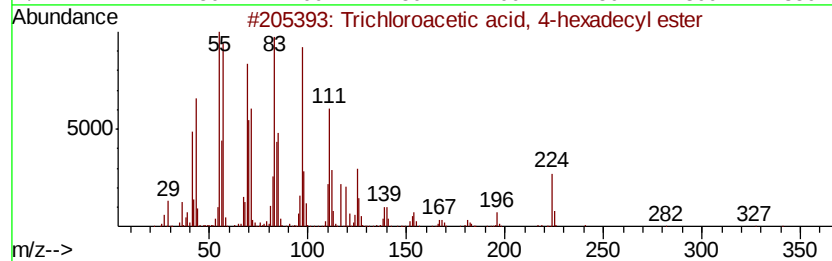
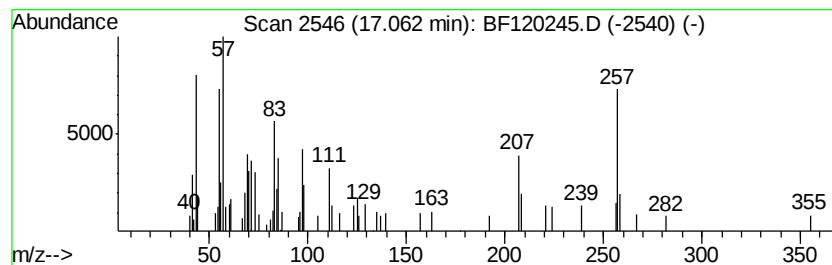
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 unknown17.06 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.06	2.03 ng	80939	Perylene-d12	15.22

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Trichloroacetic acid, 4-hexadecyl...	386	C18H33Cl3O2	1000282-92-4	46
2		Cetene	224	C16H32	000629-73-2	38
3		Hexadecyl methanesulfonate	320	C17H36O3S	020779-14-0	38
4		Ethanol, 2-(tetradecyloxy)-	258	C16H34O2	002136-70-1	38
5		Chloroacetic acid, 4-hexadecyl e...	318	C18H35ClO2	1000282-93-2	38



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF043020\
Data File : BF120245.D
Acq On : 30 Apr 2020 17:55
Operator : MA/SJ
Sample : L2451-01
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MW-1

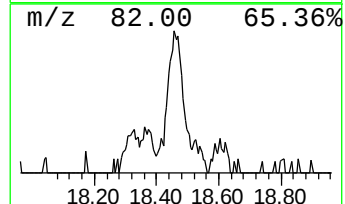
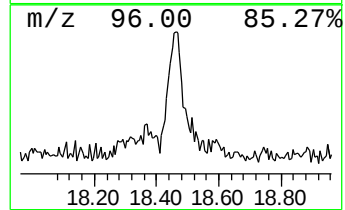
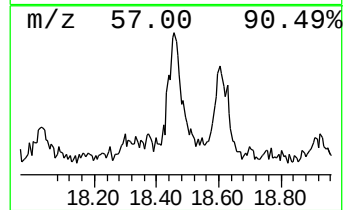
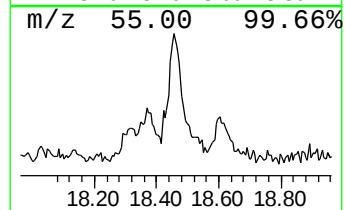
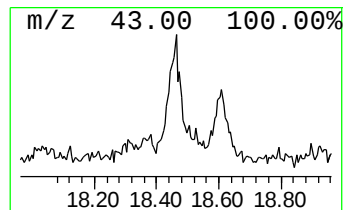
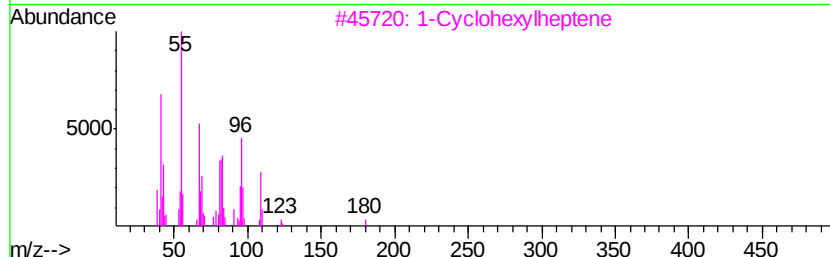
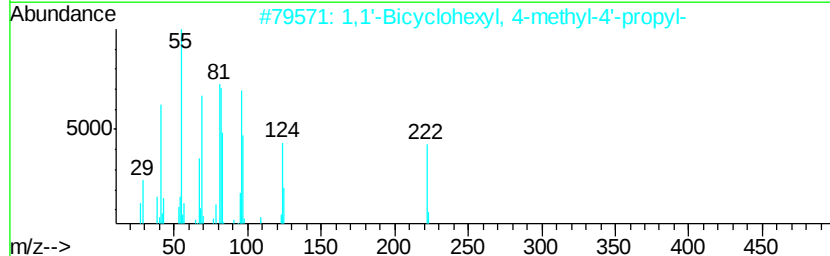
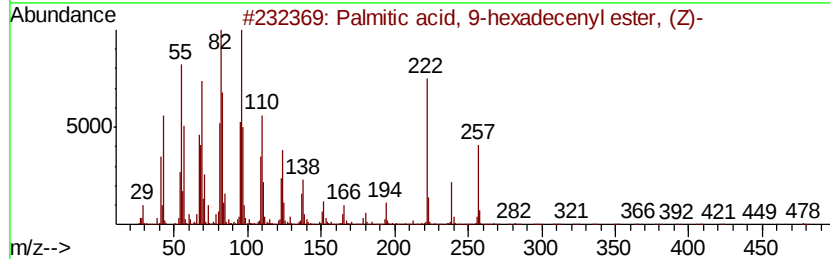
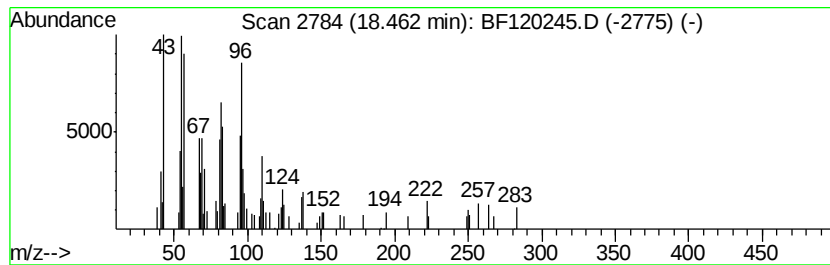
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Palmitic acid, 9-hexadecenyl... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.46	4.20 ng	167609	Perylene-d12	15.22

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Palmitic acid, 9-hexadecenyl est...	478	C32H62O2	022393-90-4	64
2		1,1'-Bicyclohexyl, 4-methyl-4'-p...	222	C16H30	092343-70-9	55
3		1-Cyclohexylheptene	180	C13H24	114614-83-4	49
4		9-Octadecen-1-ol, (Z)-	268	C18H36O	000143-28-2	43
5		11-Hexadecen-1-ol, acetate, (Z)-	282	C18H34O2	034010-21-4	38



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF043020\
 Data File : BF120245.D
 Acq On : 30 Apr 2020 17:55
 Operator : MA/SJ
 Sample : L2451-01
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MW-1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF040820.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
3-Phenylpropanol	8.17	4.2 ng		147309	2	7.93	707125	20.0
unknown8.66	8.66	9.5 ng		336993	2	7.93	707125	20.0
Phenol, 2,6-dimet...	8.85	2.1 ng		89009	3	9.69	839898	20.0
Phenol, 3,4,5-tri...	10.14	2.4 ng		99010	3	9.69	839898	20.0
Benzenemethanol, ...	10.42	2.0 ng		85899	3	9.69	839898	20.0
unknown10.92	10.92	12.4 ng		528418	4	11.17	849338	20.0
n-Hexadecanoic acid	11.72	6.4 ng		271579	4	11.17	849338	20.0
Octadecanoic acid	12.50	2.7 ng		110226	5	13.82	826651	20.0
Cyclopropanecarbo...	12.65	5.4 ng		221381	5	13.82	826651	20.0
Trifluoroacetoxy ...	13.67	5.5 ng		227072	5	13.82	826651	20.0
1,15-Hexadecadiene	16.96	2.1 ng		84452	6	15.22	797838	20.0
unknown17.06	17.06	2.0 ng		80939	6	15.22	797838	20.0
Palmitic acid, 9-...	18.46	4.2 ng		167609	6	15.22	797838	20.0