

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF101921\
 Data File : BF125897.D
 Acq On : 19 Oct 2021 13:39
 Operator : JU/CG
 Sample : M4248-01
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CHVB-0513

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-BF100721.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF125897.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.439	565	570	573	rBV	2963399	2827696	84.55%	9.493%
2	6.451	736	742	745	rBV	2695628	2773332	82.92%	9.311%
3	6.822	801	805	808	rBV	1117033	862019	25.78%	2.894%
4	7.381	895	900	903	rBV	2009690	1761799	52.68%	5.915%
5	8.004	1003	1006	1013	rBV	442728	367167	10.98%	1.233%
6	8.104	1017	1023	1026	rBV	1538179	1199562	35.87%	4.027%
7	8.980	1170	1172	1176	rBV5	47335	48223	1.44%	0.162%
8	9.063	1181	1186	1189	rVV4	41883	72111	2.16%	0.242%
9	9.127	1194	1197	1201	rVB	185460	160095	4.79%	0.537%
10	9.180	1201	1206	1209	rBV	3828219	3344391	100.00%	11.228%
11	9.269	1215	1221	1224	rBV5	171015	281958	8.43%	0.947%
12	9.410	1243	1245	1246	rBV	60807	49492	1.48%	0.166%
13	9.492	1255	1259	1262	rBV3	201912	275370	8.23%	0.924%
14	9.533	1264	1266	1270	rVB3	121070	139096	4.16%	0.467%
15	9.586	1270	1275	1277	rBV2	441271	535273	16.01%	1.797%
16	9.733	1296	1300	1302	rBV4	174150	230429	6.89%	0.774%
17	9.769	1302	1306	1309	rVV4	219245	318134	9.51%	1.068%
18	9.857	1315	1321	1325	rBV2	2051091	2132881	63.77%	7.161%
19	10.116	1363	1365	1368	rBV3	254050	228916	6.84%	0.769%
20	10.174	1373	1375	1376	rBV	205711	167809	5.02%	0.563%
21	10.269	1388	1391	1393	rBV	258441	261260	7.81%	0.877%
22	10.516	1430	1433	1439	rVB	606500	631291	18.88%	2.119%
23	10.645	1451	1455	1458	rBV	1720781	1640774	49.06%	5.508%
24	10.780	1475	1478	1485	rVB	1209974	1185617	35.45%	3.980%
25	11.245	1554	1557	1562	rVB	1524309	1452738	43.44%	4.877%
26	11.345	1571	1574	1579	rVB	1280712	1325261	39.63%	4.449%
27	11.592	1614	1616	1619	rVB2	466120	398735	11.92%	1.339%
28	12.927	1838	1843	1846	rBV	3107538	3125506	93.46%	10.493%
29	13.968	2017	2020	2025	rVB	1053074	1040276	31.11%	3.492%
30	15.080	2206	2209	2213	rVB2	39322	45955	1.37%	0.154%
31	15.421	2262	2267	2271	rBV	763363	903171	27.01%	3.032%

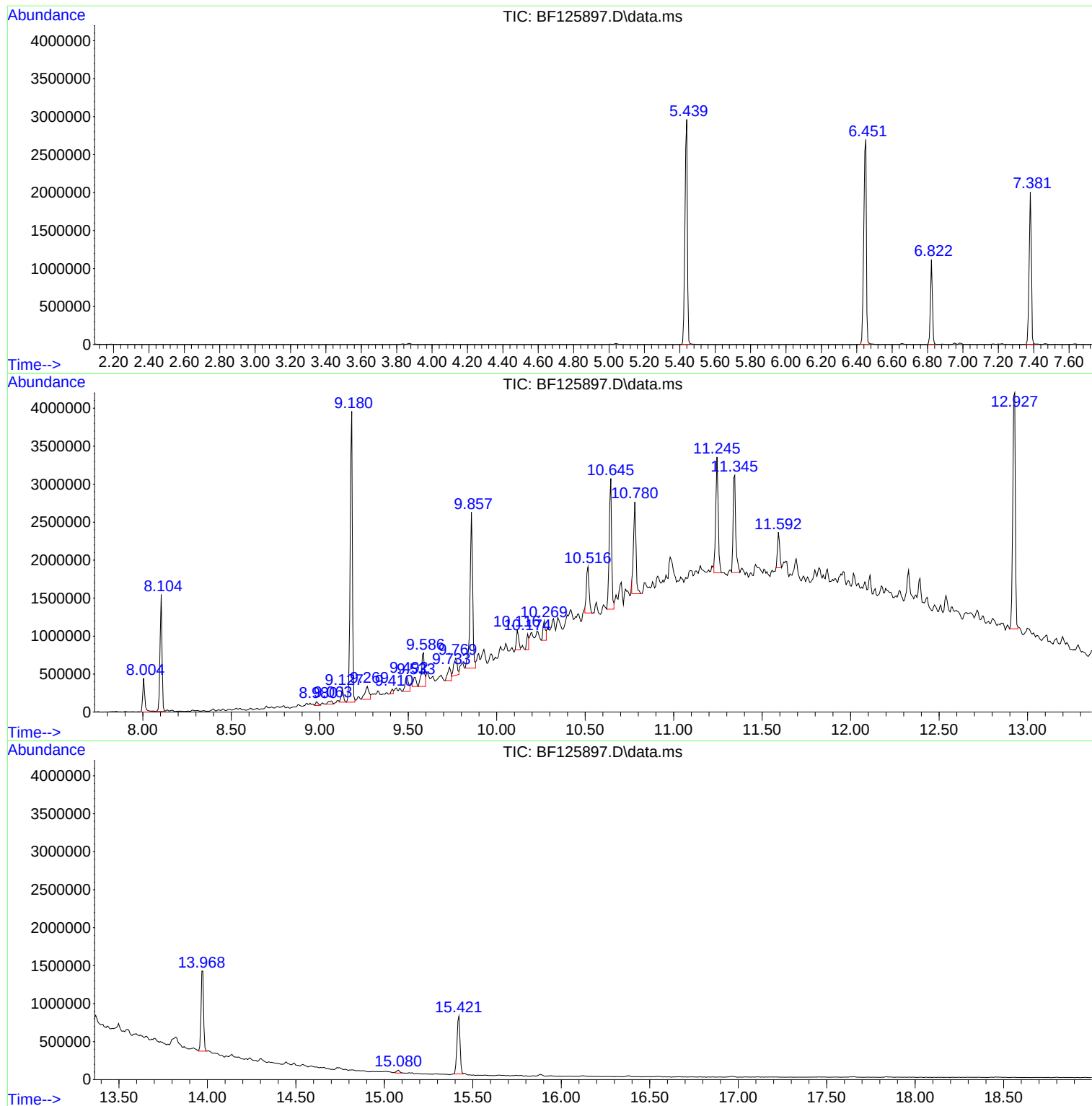
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF100721.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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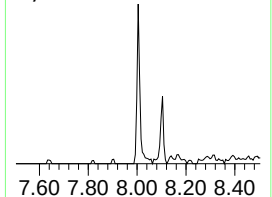
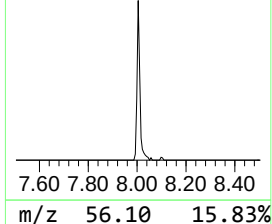
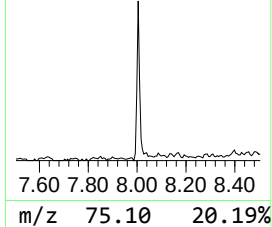
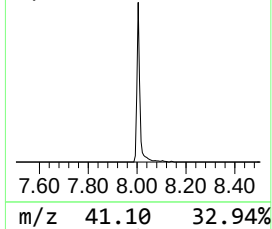
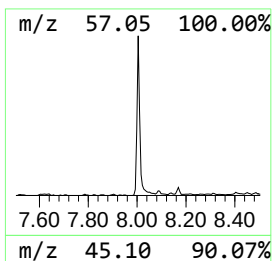
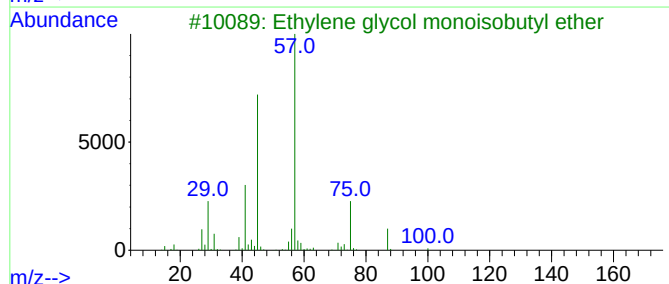
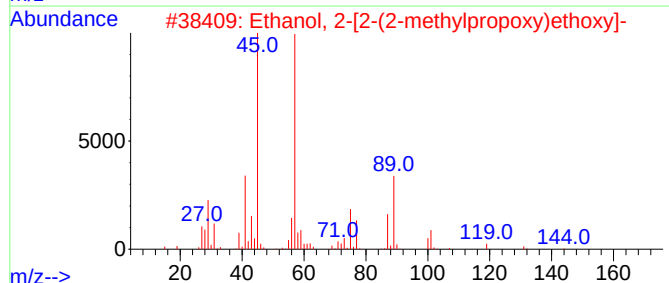
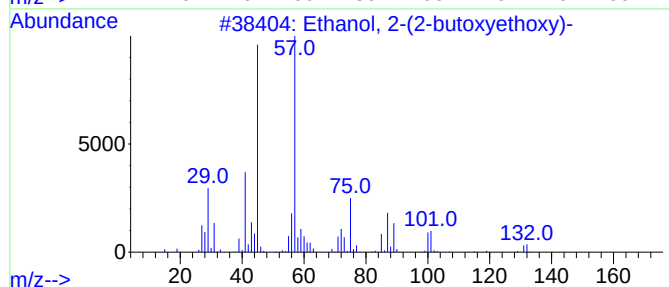
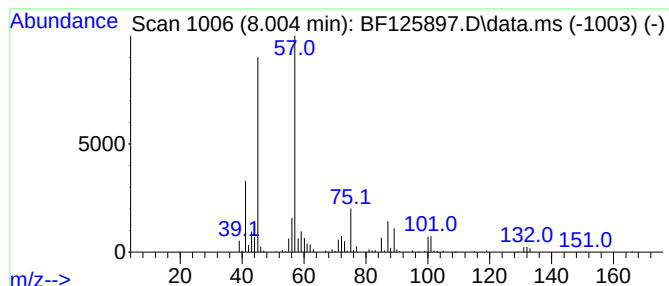
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF100721.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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Peak Number 1 Ethanol, 2-(2-butoxyethoxy)- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.004	6.12 ng	367167	Naphthalene-d8	8.104

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	91
2			Ethanol, 2-[2-(2-methylpropoxy)e...	162	C8H18O3	018912-80-6	72
3			Ethylene glycol monoisobutyl ether	118	C6H14O2	004439-24-1	64
4			Ethanol, 1-(2-butoxyethoxy)-	162	C8H18O3	054446-78-5	50
5			3,3-Dimethylbutane-2-ol	102	C6H14O	000464-07-3	50



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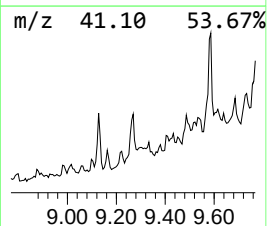
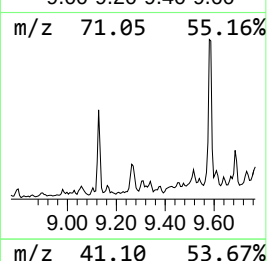
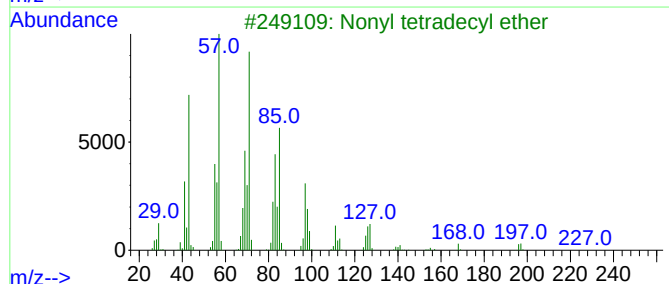
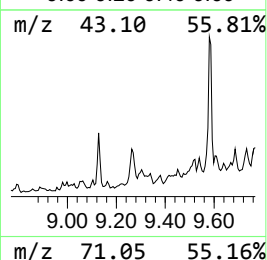
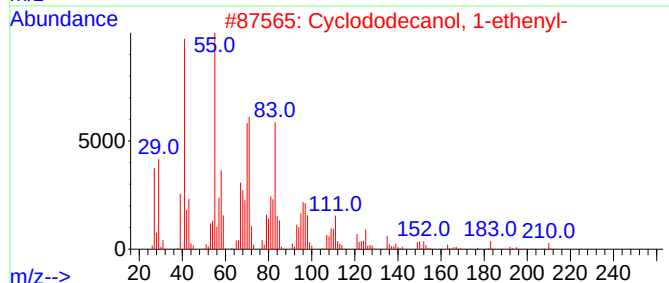
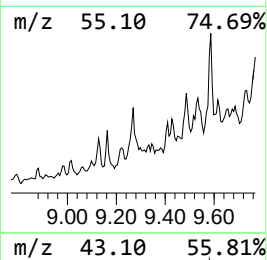
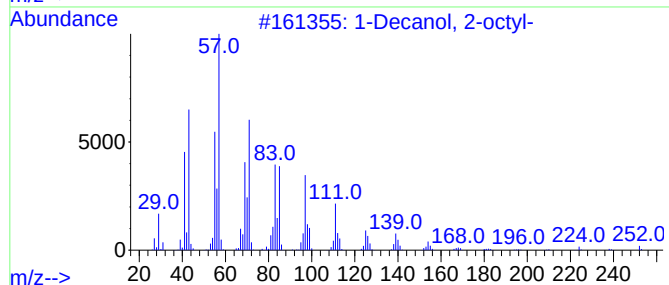
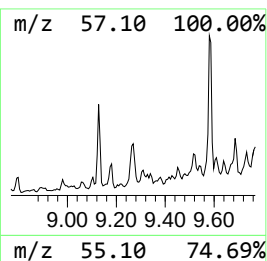
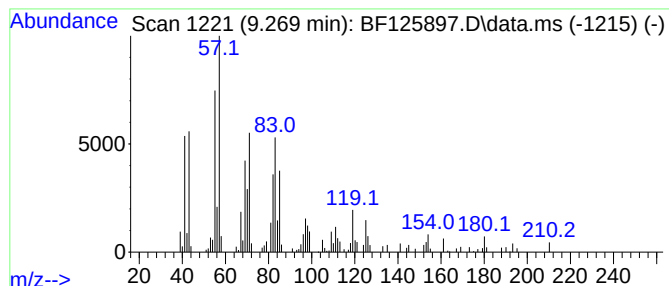
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Peak Number 2 1-Decanol, 2-octyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.269	2.64 ng	281958	Acenaphthene-d10	9.857

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Decanol, 2-octyl-			270	C18H38O	045235-48-1	64
2	Cyclododecanol, 1-ethenyl-			210	C14H26O	006244-49-1	64
3	Nonyl tetradecyl ether			340	C23H48O	1000406-37-6	52
4	Dodecyl nonyl ether			312	C21H44O	1000406-37-5	52
5	1-Dotriacontanol, acetate			509	C34H68O2	023811-94-1	52



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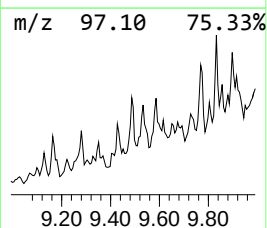
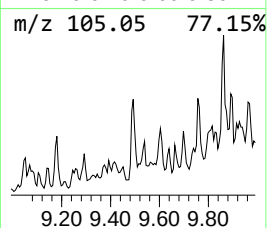
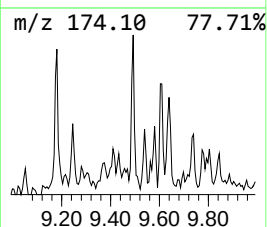
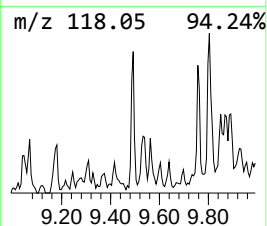
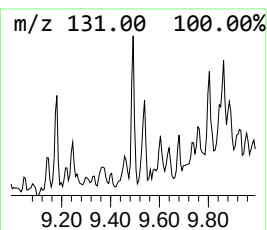
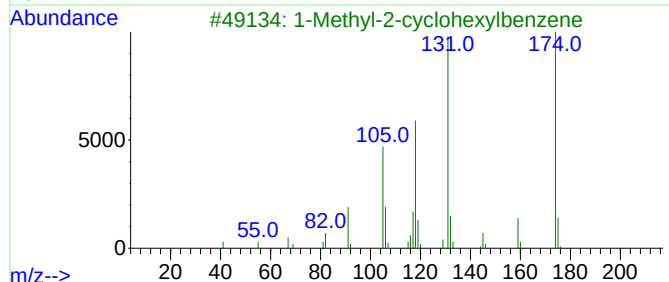
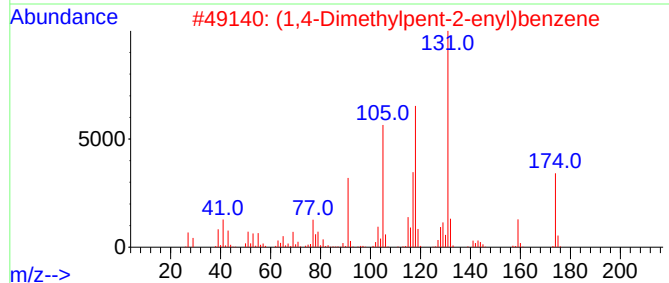
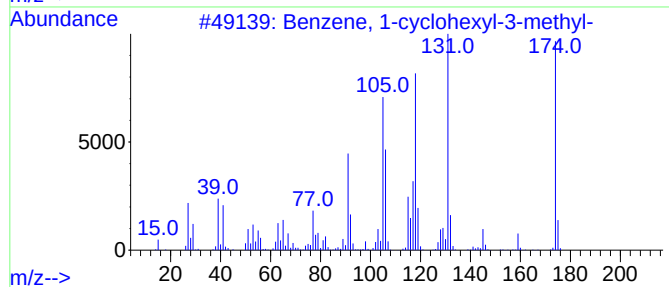
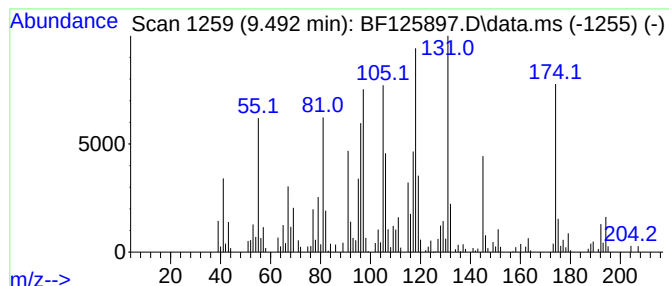
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Peak Number 3 Benzene, 1-cyclohexyl-3-met... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.492	2.58 ng	275370	Acenaphthene-d10	9.857

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-cyclohexyl-3-methyl-	174	C13H18	004575-46-6	91
2			(1,4-Dimethylpent-2-enyl)benzene	174	C13H18	1000184-97-9	55
3			1-Methyl-2-cyclohexylbenzene	174	C13H18	004501-35-3	47
4			2-Pentanone, 3-(phenylmethylene)-	174	C12H14O	003437-89-6	18
5			7-Methoxy-1-naphthol	174	C11H10O2	067247-13-6	14



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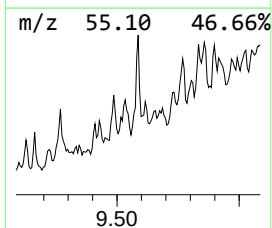
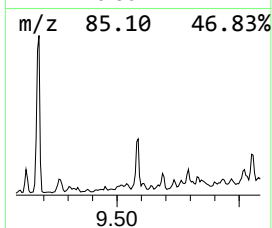
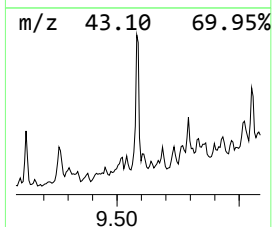
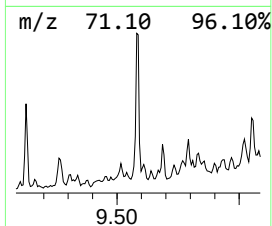
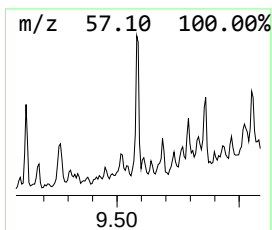
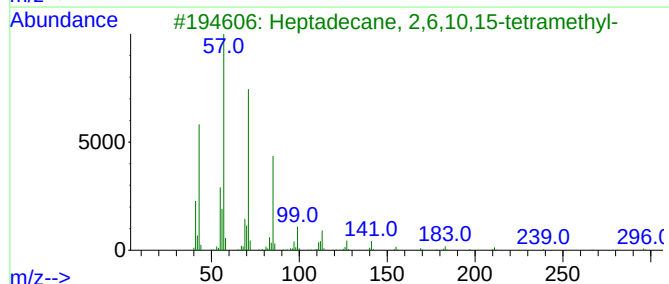
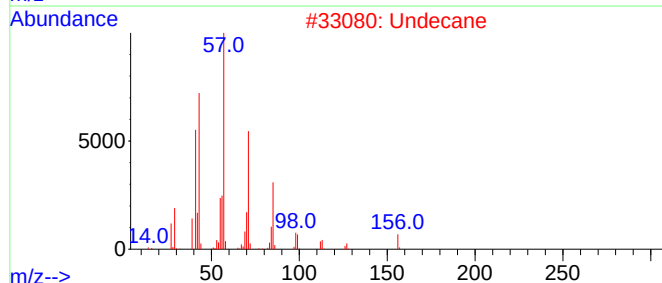
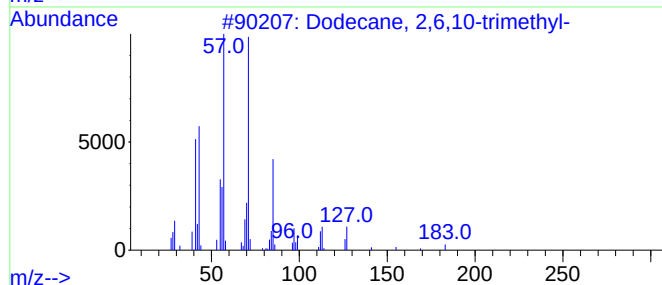
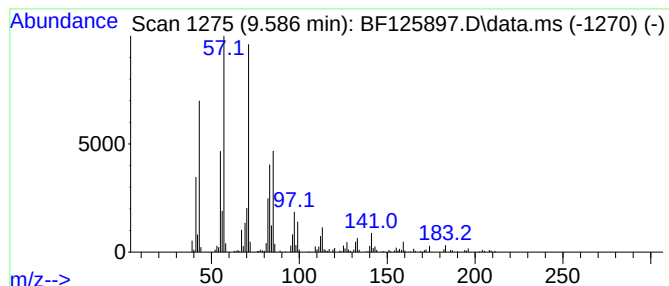
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TIC Library : C:\Database\NIST20.L
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Peak Number 4 Dodecane, 2,6,10-trimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.586	5.02 ng	535273	Acenaphthene-d10	9.857

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dodecane, 2,6,10-trimethyl-	212	C15H32	003891-98-3	64
2			Undecane	156	C11H24	001120-21-4	55
3			Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	52
4			Dodecane, 1-iodo-	296	C12H25I	004292-19-7	52
5			Hexadecane	226	C16H34	000544-76-3	47



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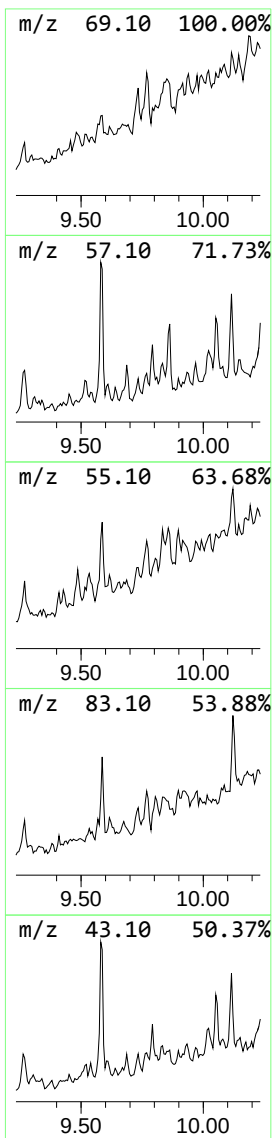
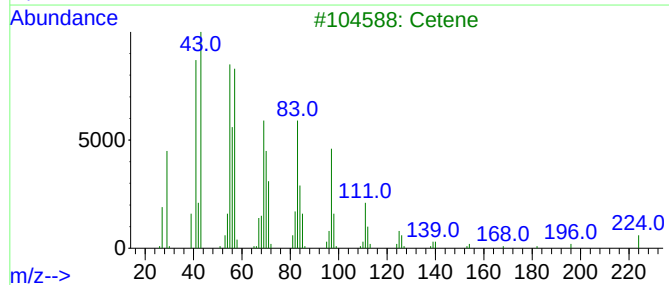
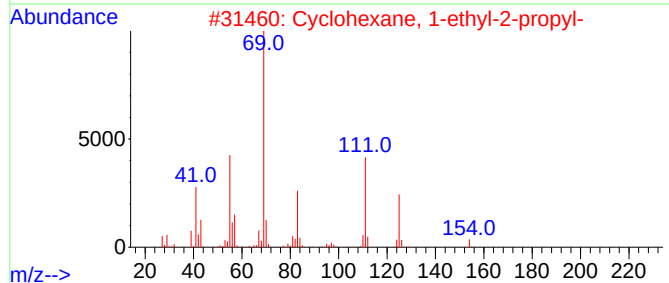
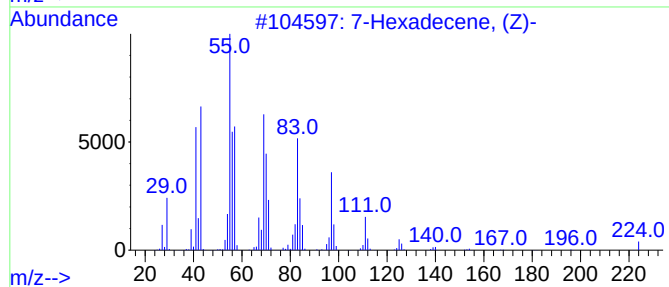
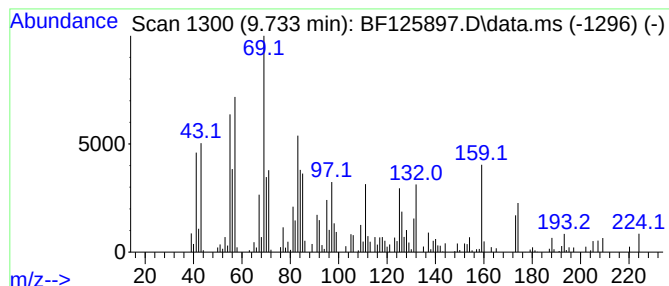
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Peak Number 5 unknown9.733 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.733	2.16 ng	230429	Acenaphthene-d10	9.857

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	7-Hexadecene, (Z)-			224	C16H32	035507-09-6	38
2	Cyclohexane, 1-ethyl-2-propyl-			154	C11H22	062238-33-9	30
3	Cetene			224	C16H32	000629-73-2	25
4	1-Hexene, 3,3-dimethyl-			112	C8H16	003404-77-1	18
5	Hexatriacontyl pentafluoropropio...			669	C39H73F5O2	1000351-89-0	11



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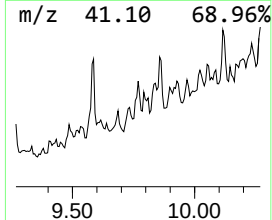
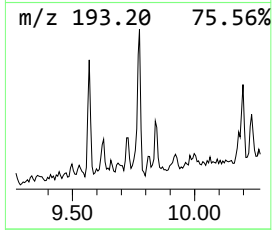
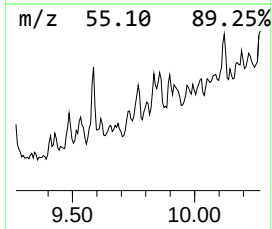
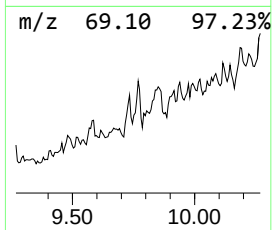
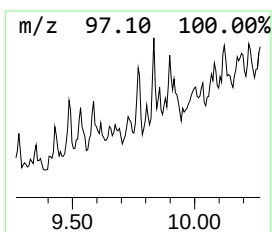
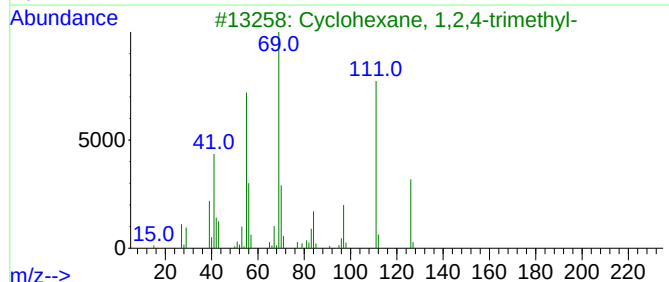
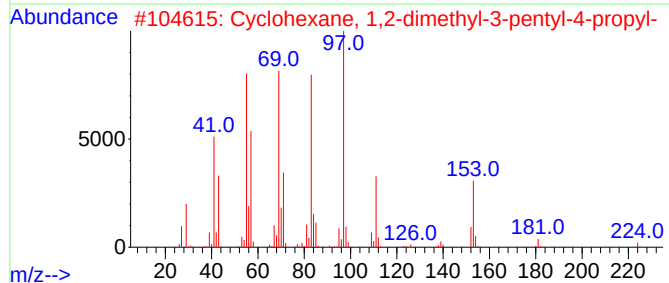
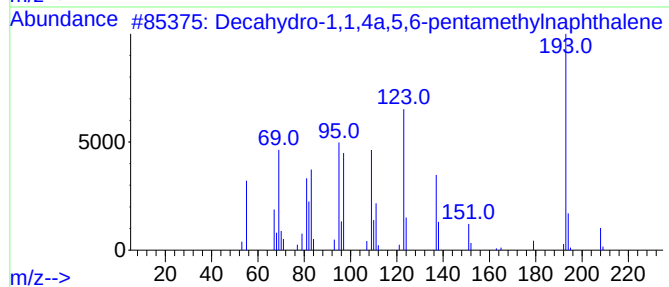
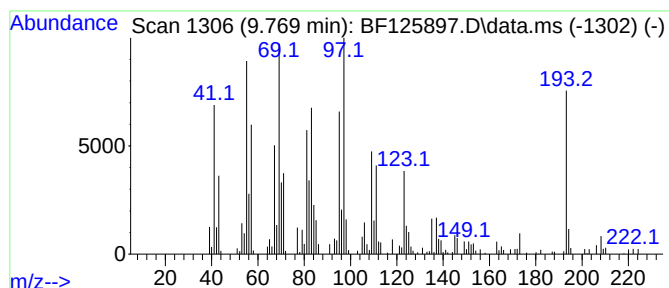
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Decahydro-1,1,4a,5,6-pentam... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.	
9.769	2.98 ng	318134	Acenaphthene-d10	9.857	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Decahydro-1,1,4a,5,6-pentamethyl...	208	C15H28	080655-44-3	50
2	Cyclohexane, 1,2-dimethyl-3-pent...	224	C16H32	062376-17-4	46
3	Cyclohexane, 1,2,4-trimethyl-	126	C9H18	002234-75-5	30
4	6-Bromo-2,2-dimethylcyclohexanone	204	C8H13BrO	021690-26-6	22
5	1,7-Dimethyl-4-(1-methylethyl)cv...	210	C15H30	000645-10-3	22



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF101921\
Data File : BF125897.D
Acq On : 19 Oct 2021 13:39
Operator : JU/CG
Sample : M4248-01
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
CHVB-0513

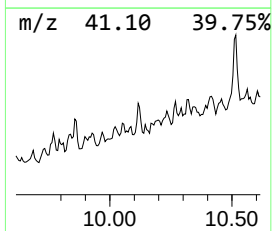
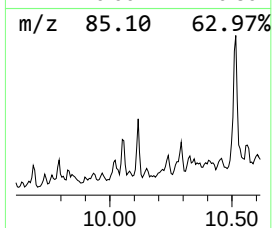
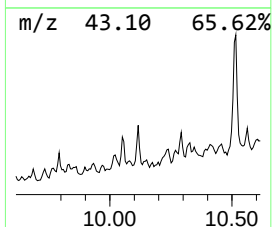
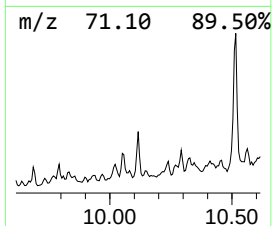
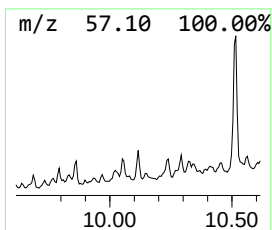
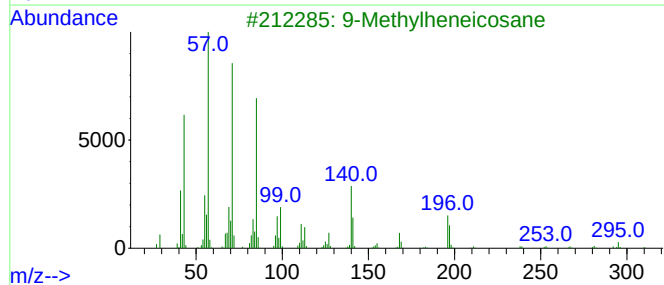
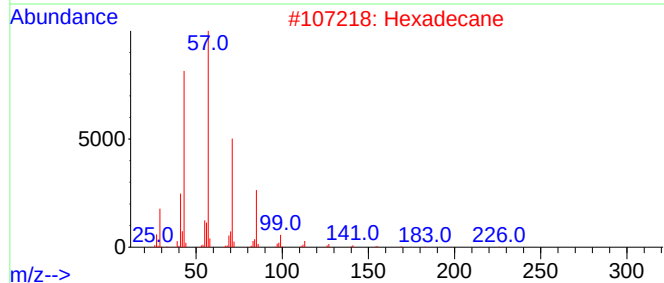
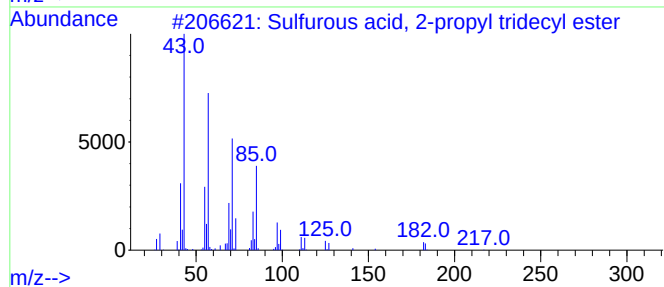
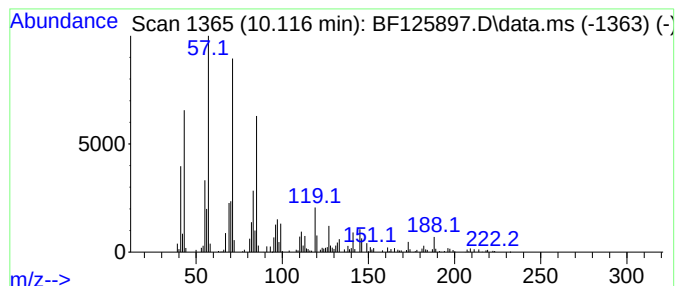
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Sulfurous acid, 2-propyl tr... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.116	2.15 ng	228916	Acenaphthene-d10	9.857

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Sulfurous acid, 2-propyl tridecy...	306	C16H34O3S	1000309-12-4	59
2			Hexadecane	226	C16H34	000544-76-3	50
3			9-Methylheneicosane	310	C22H46	070475-51-3	47
4			Heptacosane	380	C27H56	000593-49-7	47
5			Succinic acid, 2-chloro-6-fluoro...	330	C15H16ClF05	1000390-72-1	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF101921\
Data File : BF125897.D
Acq On : 19 Oct 2021 13:39
Operator : JU/CG
Sample : M4248-01
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
CHVB-0513

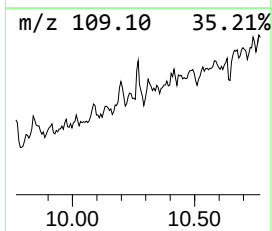
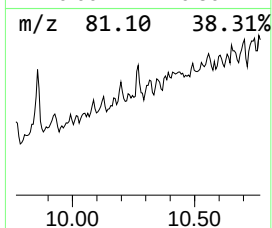
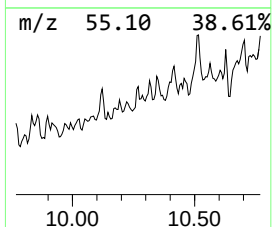
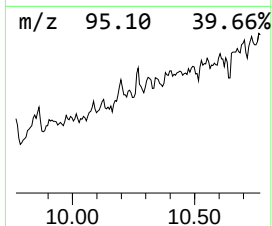
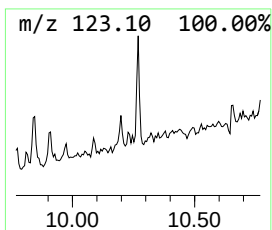
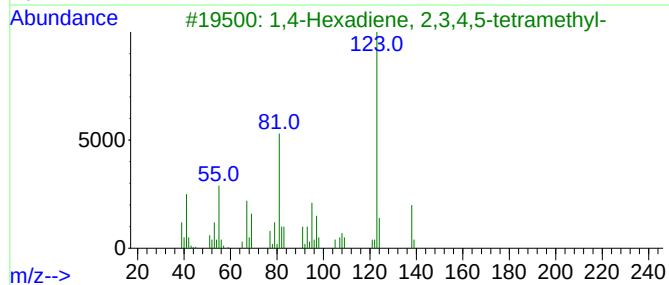
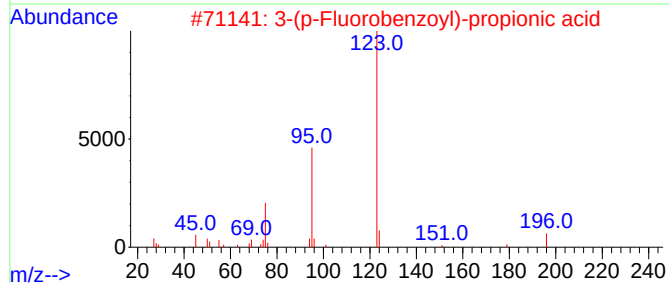
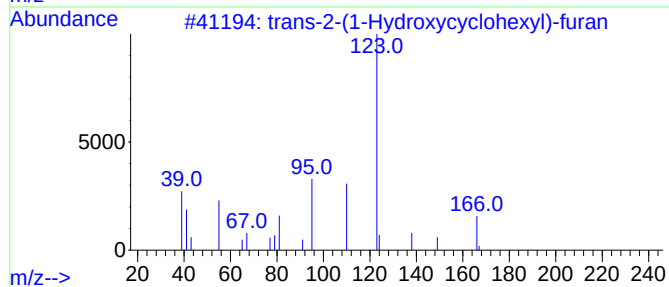
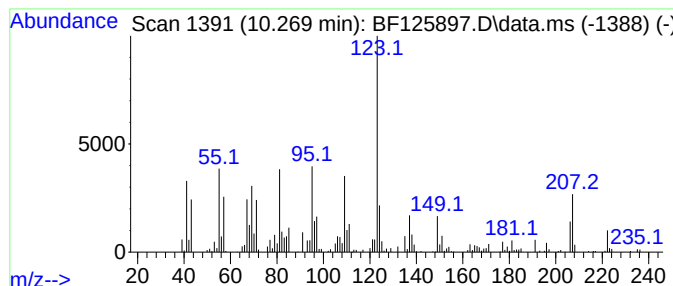
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 unknown10.269 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.269	2.45 ng	261260	Acenaphthene-d10	9.857

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			trans-2-(1-Hydroxycyclohexyl)-furan	166	C10H14O2	115754-87-5	47
2			3-(p-Fluorobenzoyl)-propionic acid	196	C10H9FO3	000366-77-8	47
3			1,4-Hexadiene, 2,3,4,5-tetramethyl-	138	C10H18	051504-54-2	47
4			2-(4-Methoxyphenoxy)propanoic acid	196	C10H12O4	013794-15-5	47
5			3-Fluorobenzoic acid, 2-ethylcyc...	250	C15H19FO2	1000279-04-9	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF101921\
Data File : BF125897.D
Acq On : 19 Oct 2021 13:39
Operator : JU/CG
Sample : M4248-01
Misc :
ALS Vial : 7 Sample Multiplier: 1

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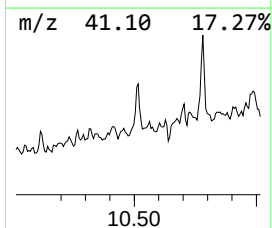
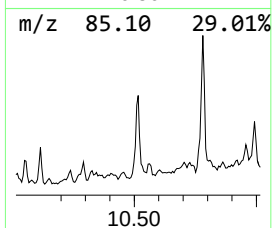
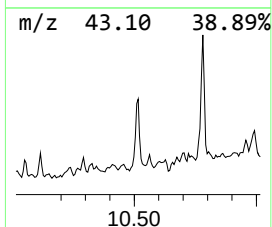
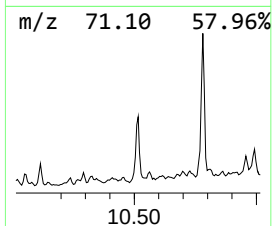
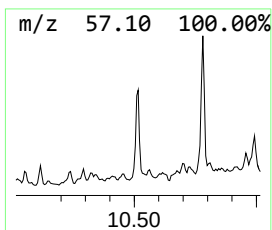
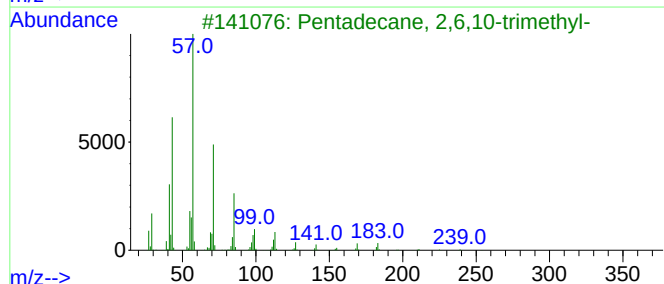
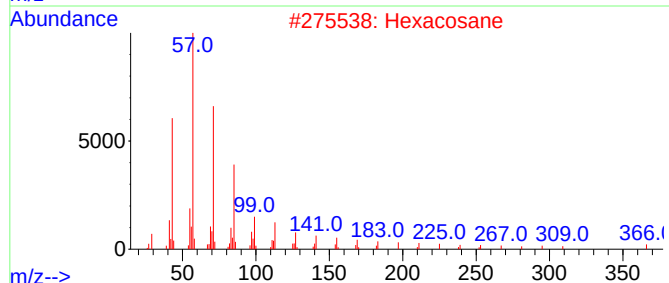
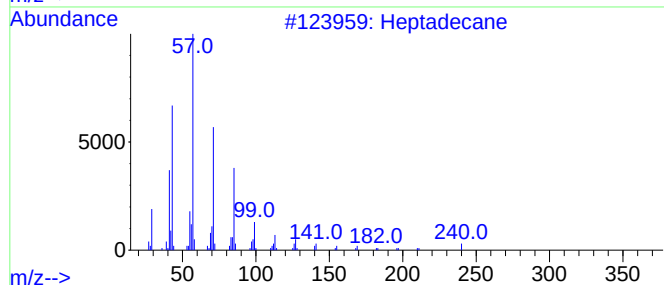
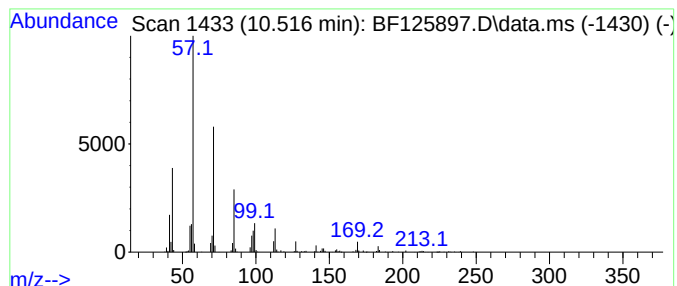
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Heptadecane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.516	5.92 ng	631291	Acenaphthene-d10	9.857

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptadecane	240	C17H36	000629-78-7	72
2			Hexacosane	366	C26H54	000630-01-3	72
3			Pentadecane, 2,6,10-trimethyl-	254	C18H38	003892-00-0	64
4			Tetratetracontane	619	C44H90	007098-22-8	64
5			Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF101921\
Data File : BF125897.D
Acq On : 19 Oct 2021 13:39
Operator : JU/CG
Sample : M4248-01
Misc :
ALS Vial : 7 Sample Multiplier: 1

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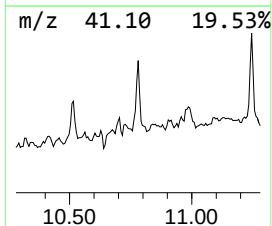
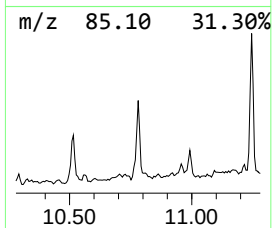
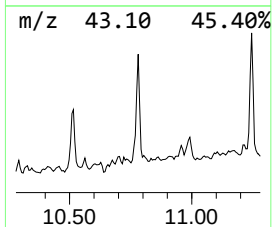
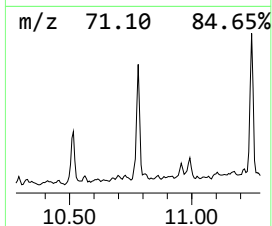
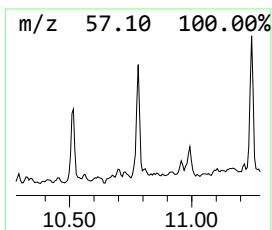
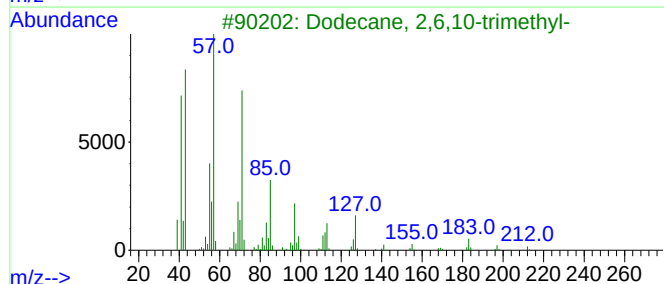
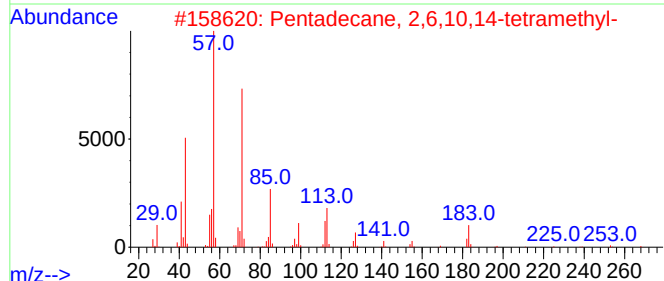
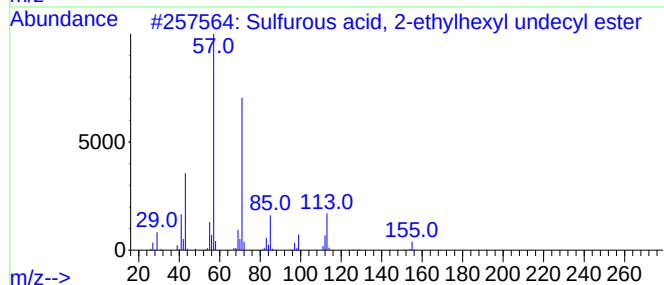
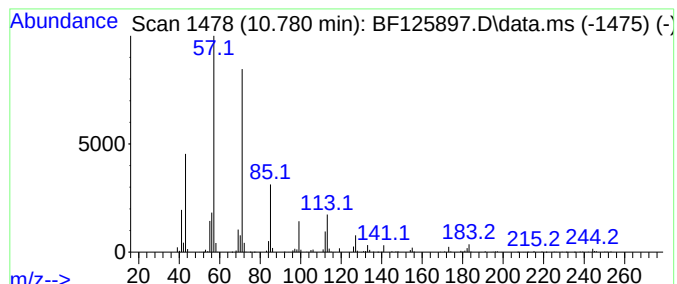
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Sulfurous acid, 2-ethylhexy... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.780	17.89 ng	1185620	Phenanthrene-d10	11.339

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Sulfurous acid, 2-ethylhexyl und...	348	C19H40O3S	1000309-19-4	80
2			Pentadecane, 2,6,10,14-tetramethyl-	268	C19H40	001921-70-6	78
3			Dodecane, 2,6,10-trimethyl-	212	C15H32	003891-98-3	72
4			Sulfurous acid, 2-ethylhexyl tri...	376	C21H44O3S	1000309-19-6	72
5			Undecane, 6,6-dimethyl-	184	C13H28	017312-76-4	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF101921\
Data File : BF125897.D
Acq On : 19 Oct 2021 13:39
Operator : JU/CG
Sample : M4248-01
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
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ClientSampleId :
CHVB-0513

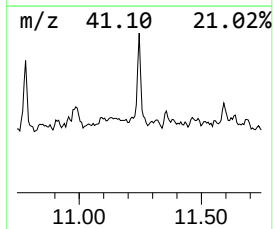
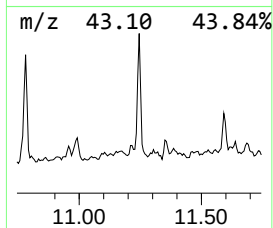
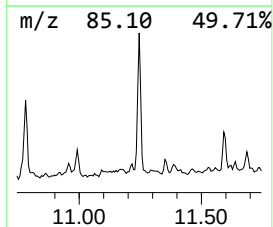
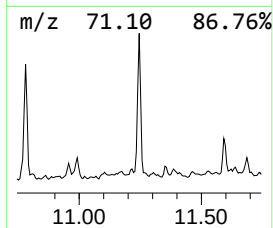
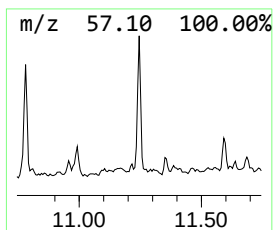
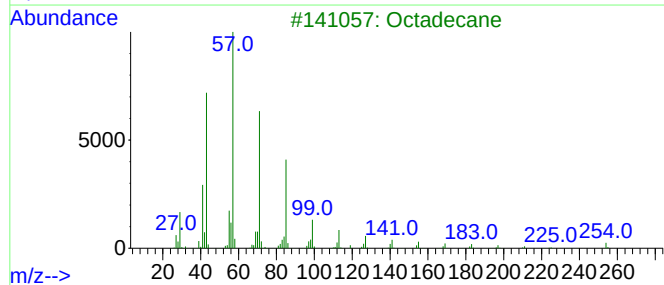
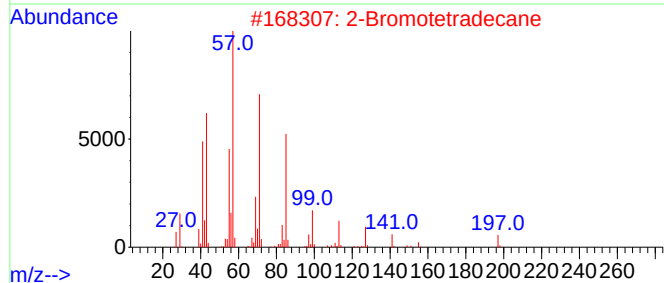
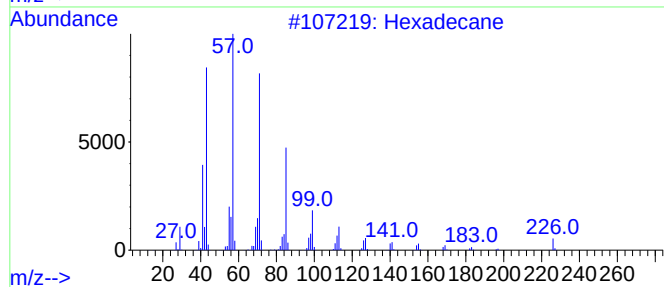
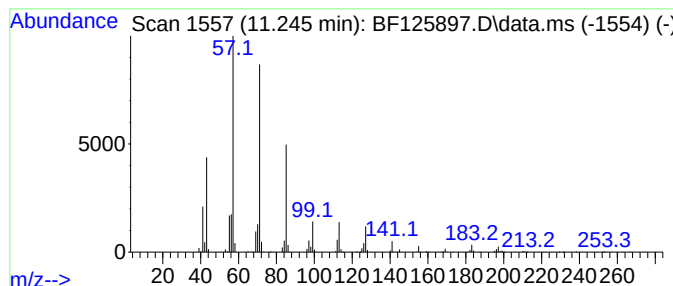
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Hexadecane Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.245	21.92 ng	1452740	Phenanthrene-d10	11.339

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexadecane	226	C16H34	000544-76-3	91
2			2-Bromotetradecane	276	C14H29Br	074036-95-6	90
3			Octadecane	254	C18H38	000593-45-3	90
4			Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	87
5			Tritetracontane	605	C43H88	007098-21-7	78



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF101921\
Data File : BF125897.D
Acq On : 19 Oct 2021 13:39
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Sample : M4248-01
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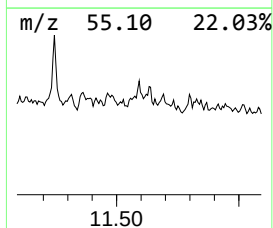
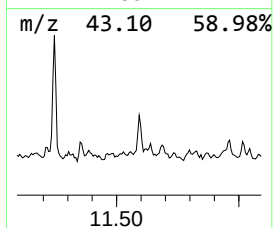
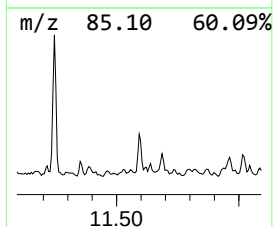
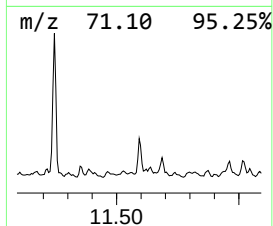
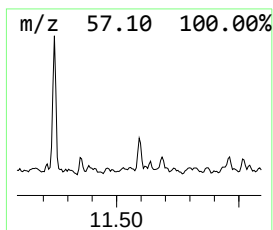
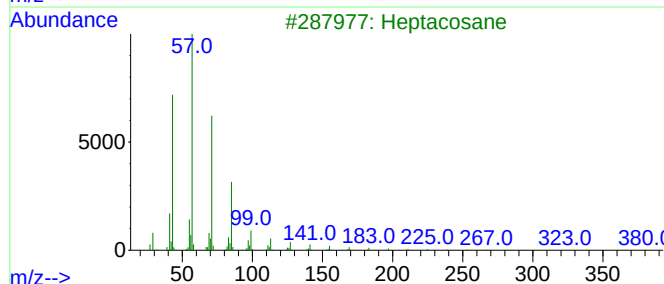
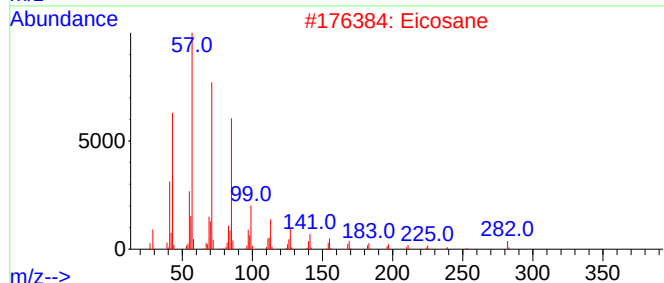
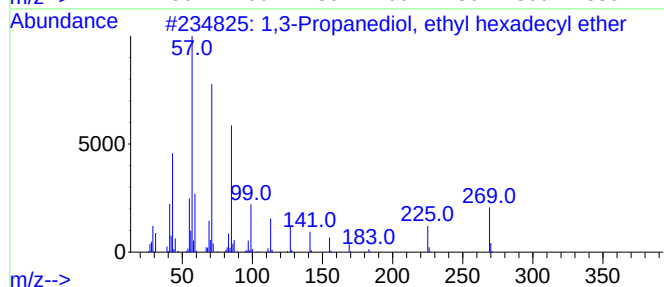
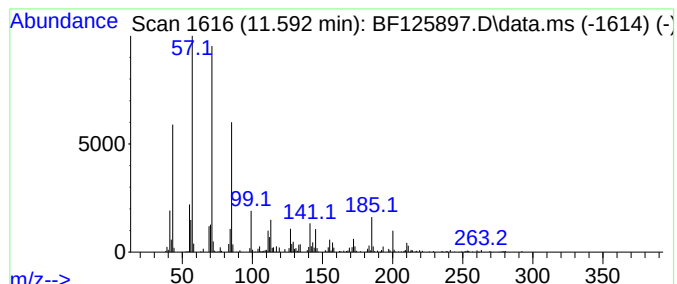
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 1,3-Propanediol, ethyl hexa... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.592	6.02 ng	398735	Phenanthrene-d10	11.339

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		1,3-Propanediol, ethyl hexadecyl...	328	C21H44O2	1000406-35-3	74
2		Eicosane	282	C20H42	000112-95-8	72
3		Heptacosane	380	C27H56	000593-49-7	72
4		Octacosane	394	C28H58	000630-02-4	72
5		Hexacosane	366	C26H54	000630-01-3	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF101921\
Data File : BF125897.D
Acq On : 19 Oct 2021 13:39
Operator : JU/CG
Sample : M4248-01
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
CHVB-0513

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF100721.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
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1-Decanol, 2-oc...	9.269	2.6	ng	281958	3	9.857	2132880	20.0
Benzene, 1-cycl...	9.492	2.6	ng	275370	3	9.857	2132880	20.0
Dodecane, 2,6,1...	9.586	5.0	ng	535273	3	9.857	2132880	20.0
unknown9.733	9.733	2.2	ng	230429	3	9.857	2132880	20.0
Decahydro-1,1,4...	9.769	3.0	ng	318134	3	9.857	2132880	20.0
Sulfurous acid,...	10.116	2.1	ng	228916	3	9.857	2132880	20.0
unknown10.269	10.269	2.5	ng	261260	3	9.857	2132880	20.0
Heptadecane	10.516	5.9	ng	631291	3	9.857	2132880	20.0
Sulfurous acid,...	10.780	17.9	ng	1185620	4	11.339	1325260	20.0
Hexadecane	11.245	21.9	ng	1452740	4	11.339	1325260	20.0
1,3-Propanediol...	11.592	6.0	ng	398735	4	11.339	1325260	20.0