

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF052920\
 Data File : BF120446.D
 Acq On : 29 May 2020 18:28
 Operator : MA/SJ
 Sample : L2773-20
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 NM73-COMP-2020-0-6

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-BF052820.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	1.934	4	8	10	rVV	543536	560751	10.83%	1.175%
2	1.958	10	12	33	rVB	1758641	2623017	50.67%	5.497%
3	2.110	33	38	55	rVB	1103715	1499541	28.97%	3.143%
4	5.451	601	606	609	rBV	3920686	3647410	70.46%	7.644%
5	6.463	772	778	781	rBV	3851693	3915634	75.64%	8.206%
6	6.663	809	812	819	rBV	89965	82041	1.58%	0.172%
7	6.834	838	841	845	rBV	1837665	1594274	30.80%	3.341%
8	6.992	864	868	871	rBV	4261376	3578652	69.13%	7.500%
9	7.398	931	937	940	rBV	3025817	2772904	53.56%	5.811%
10	8.116	1055	1059	1063	rBV	2427745	2047238	39.55%	4.290%
11	9.192	1237	1242	1246	rBV	5362448	5176924	100.00%	10.850%
12	9.586	1305	1309	1312	rBV	932769	712412	13.76%	1.493%
13	9.875	1353	1358	1361	rBV	2521831	2216975	42.82%	4.646%
14	10.657	1487	1491	1494	rBV	3158134	2762857	53.37%	5.790%
15	11.357	1605	1610	1612	rBV	2309144	2001562	38.66%	4.195%
16	11.380	1612	1614	1617	rVV	273248	217805	4.21%	0.456%
17	11.427	1617	1622	1626	rVB2	69911	68034	1.31%	0.143%
18	11.810	1683	1687	1691	rBV2	37559	53333	1.03%	0.112%
19	11.880	1697	1699	1703	rBV	146002	135723	2.62%	0.284%
20	11.969	1711	1714	1716	rBV2	55258	54249	1.05%	0.114%
21	12.563	1810	1815	1819	rBV	477302	451834	8.73%	0.947%
22	12.792	1848	1854	1857	rBV	402623	396028	7.65%	0.830%
23	12.939	1875	1879	1883	rVB	4340907	3960718	76.51%	8.301%
24	13.186	1913	1921	1924	rBV2	52304	92814	1.79%	0.195%
25	13.633	1993	1997	2001	rVB2	40501	51782	1.00%	0.109%
26	13.739	2010	2015	2018	rBV2	37745	53981	1.04%	0.113%
27	13.833	2028	2031	2032	rBV2	54594	57029	1.10%	0.120%
28	13.857	2032	2035	2043	rVB2	351921	516398	9.97%	1.082%
29	13.992	2050	2058	2060	rBV	1763815	1743850	33.69%	3.655%
30	14.016	2060	2062	2065	rVB	188612	145455	2.81%	0.305%
31	14.468	2136	2139	2141	rBV2	118745	104848	2.03%	0.220%
32	14.498	2141	2144	2152	rVB4	118990	192081	3.71%	0.403%
33	14.774	2186	2191	2195	rBV	75615	88544	1.71%	0.186%
34	14.916	2212	2215	2222	rBV4	66238	75970	1.47%	0.159%

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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 >
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 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	15.021	2228	2233	2236	rBV	176349	278976	5.39%	0.585%
36	15.110	2243	2248	2252	rBV2	195855	256749	4.96%	0.538%
37	15.157	2252	2256	2264	rVB4	79027	164035	3.17%	0.344%
38	15.315	2280	2283	2289	rVB	114619	137313	2.65%	0.288%
39	15.380	2289	2294	2297	rBV	138387	172510	3.33%	0.362%
40	15.445	2300	2305	2309	rVV	1298568	1542455	29.79%	3.233%
41	15.486	2309	2312	2317	rVB3	99296	141931	2.74%	0.297%
42	15.686	2342	2346	2349	rVB3	45871	59999	1.16%	0.126%
43	15.915	2380	2385	2392	rBV	206829	317121	6.13%	0.665%
44	15.992	2393	2398	2409	rVV3	84783	218642	4.22%	0.458%
45	16.345	2453	2458	2462	rBV7	32384	63291	1.22%	0.133%
46	16.415	2465	2470	2479	rVB6	53271	102954	1.99%	0.216%
47	16.704	2514	2519	2527	rVB6	60366	121122	2.34%	0.254%
48	16.874	2543	2548	2554	rBV2	58242	109601	2.12%	0.230%
49	17.304	2617	2621	2628	rVB	49067	86186	1.66%	0.181%
50	17.509	2649	2656	2670	rVB9	76435	290072	5.60%	0.608%

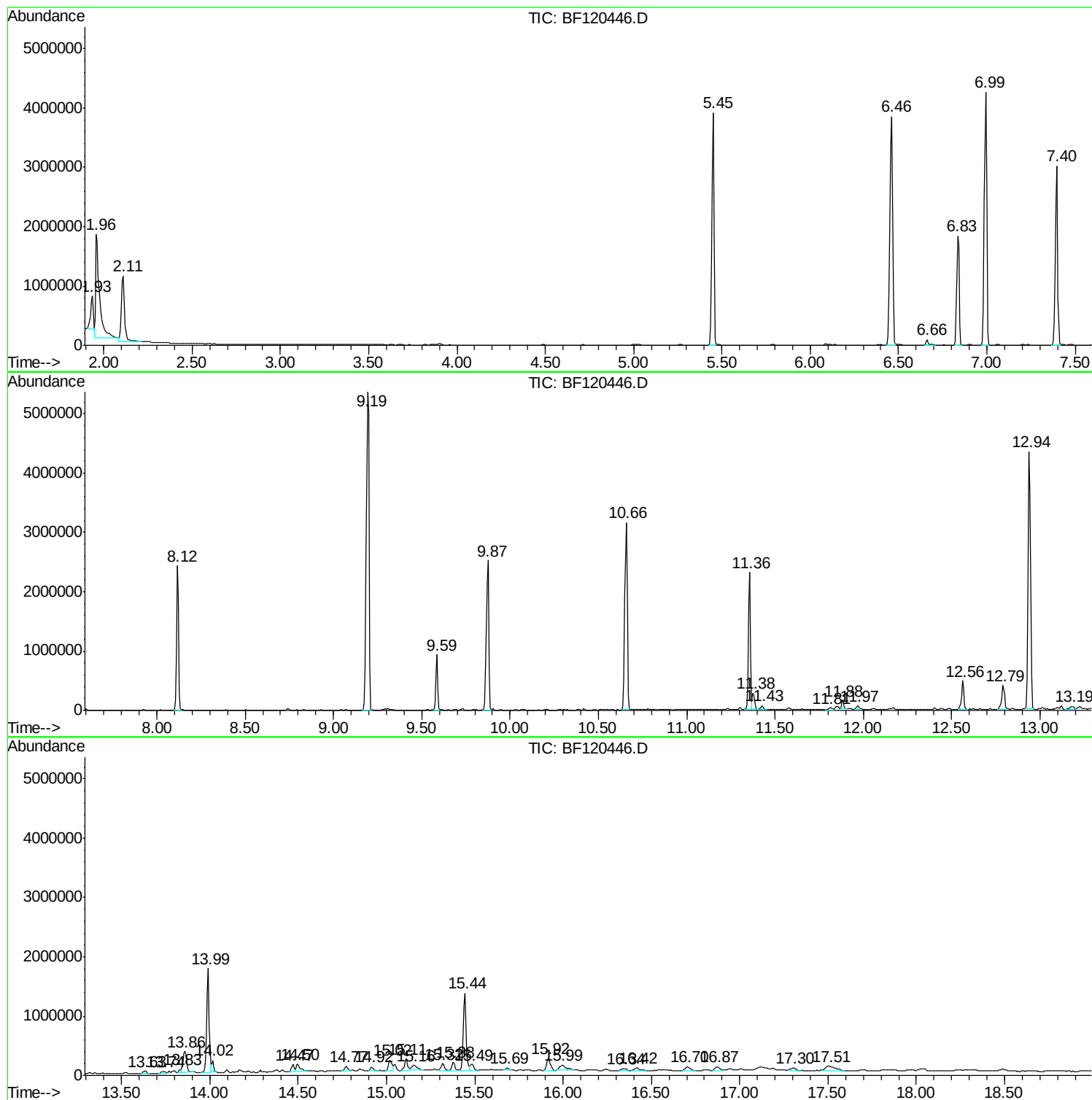
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF052820.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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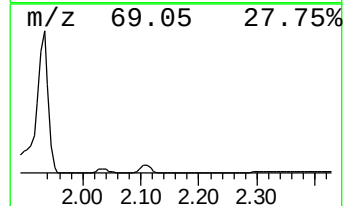
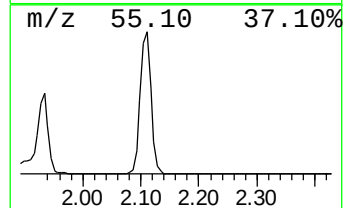
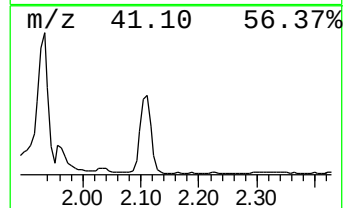
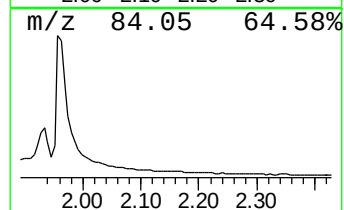
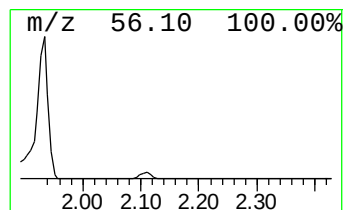
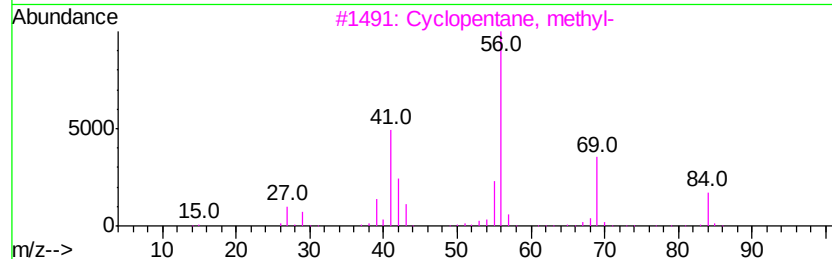
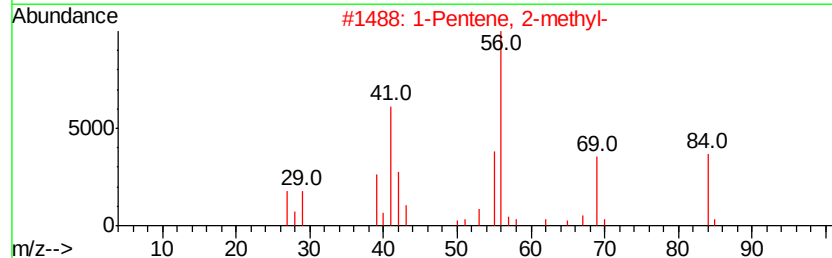
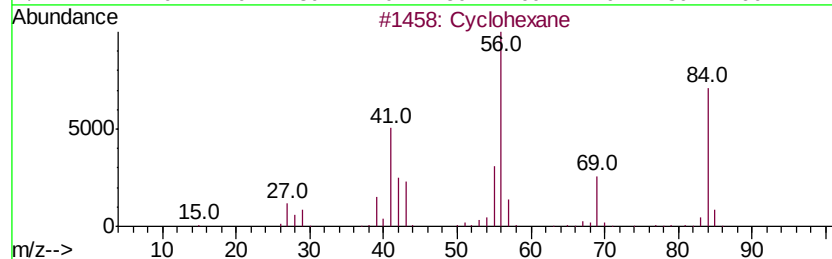
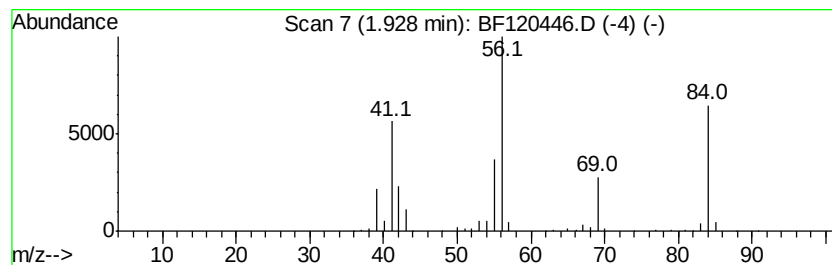
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF052820.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Cyclohexane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.93	7.03 ng	560751	1,4-Dichlorobenzene-d4	6.83

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexane	84	C6H12	000110-82-7	91
2		1-Pentene, 2-methyl-	84	C6H12	000763-29-1	64
3		Cyclopentane, methyl-	84	C6H12	000096-37-7	64
4		Cyclobutane	56	C4H8	000287-23-0	46
5		1-Hexene	84	C6H12	000592-41-6	45



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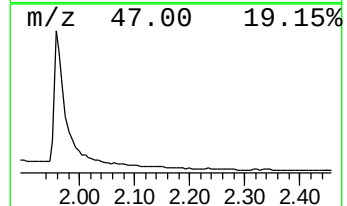
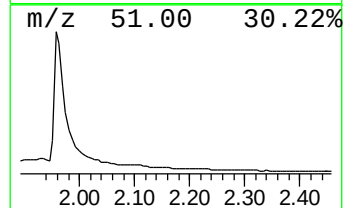
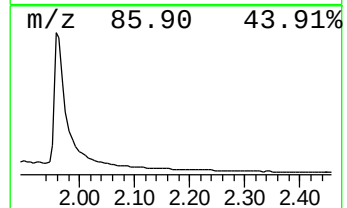
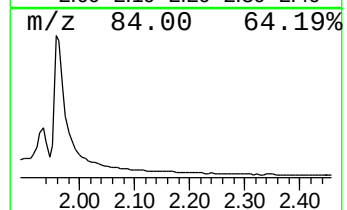
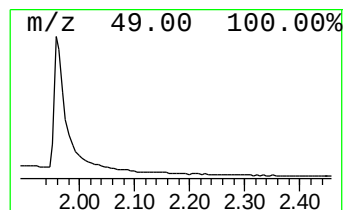
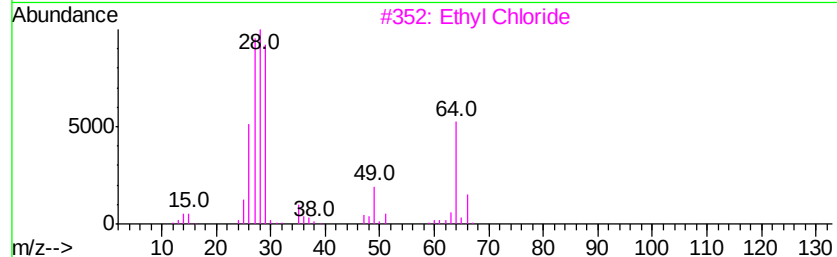
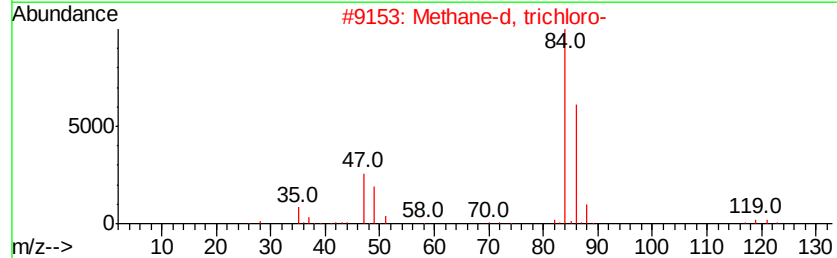
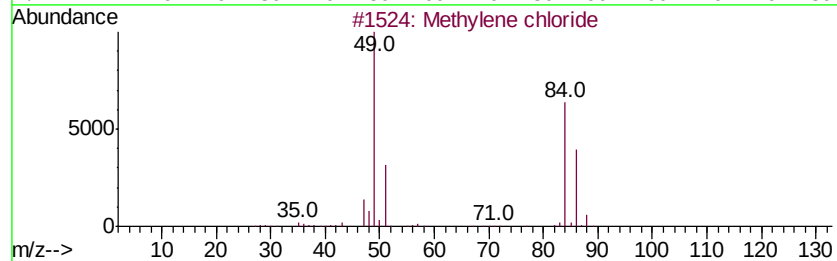
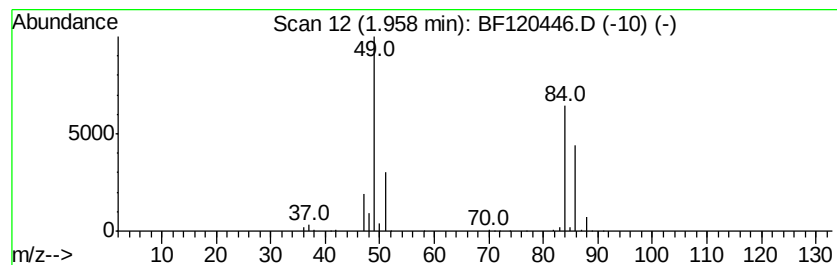
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Methylene chloride Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.96	32.91 ng	2623020	1,4-Dichlorobenzene-d4	6.83

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Methylene chloride	84	CH2Cl2	000075-09-2	95
2		Methane-d, trichloro-	119	CDCl3	000865-49-6	10
3		Ethyl Chloride	64	C2H5Cl	000075-00-3	4
4		Acetic acid, chloro-, ethyl ester	122	C4H7ClO2	000105-39-5	2
5		2-Chloroethanol	80	C2H5ClO	000107-07-3	1



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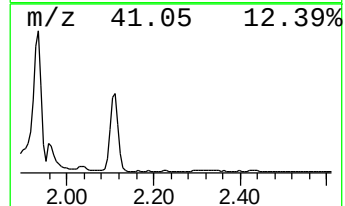
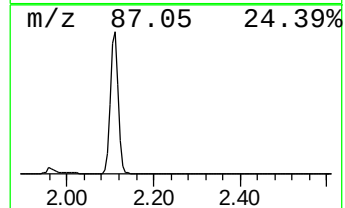
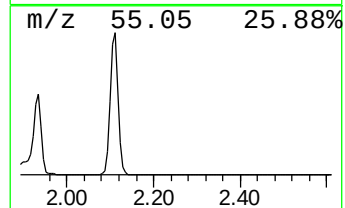
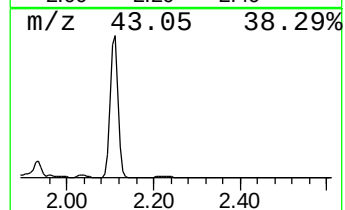
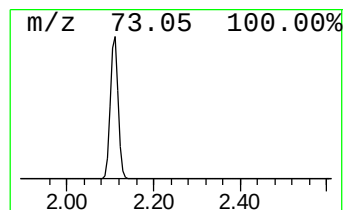
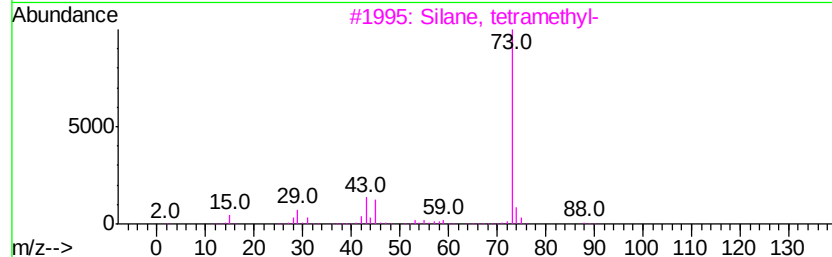
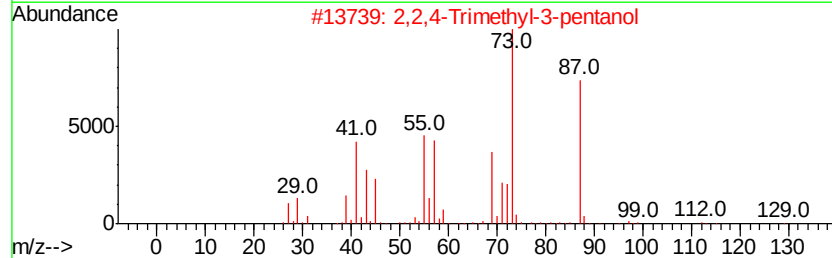
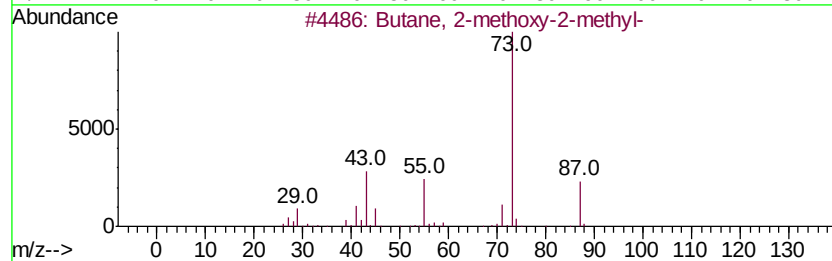
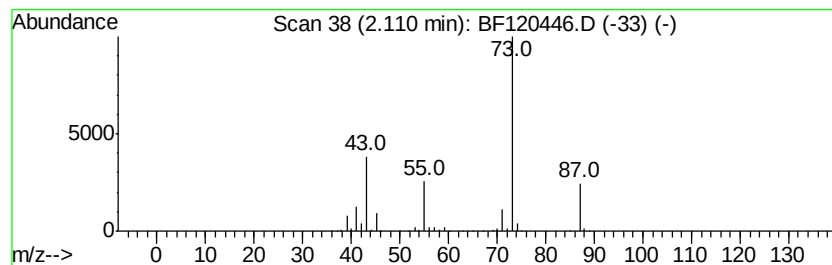
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Butane, 2-methoxy-2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.11	18.81 ng	1499540	1,4-Dichlorobenzene-d4	6.83

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2		2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39
3		Silane, tetramethyl-	88	C4H12Si	000075-76-3	35
4		1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	25
5		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	12



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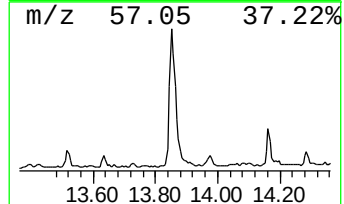
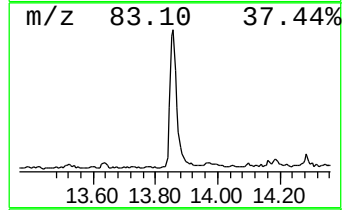
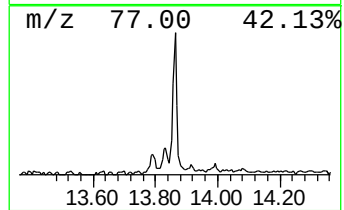
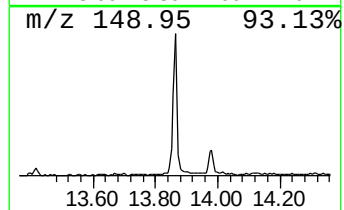
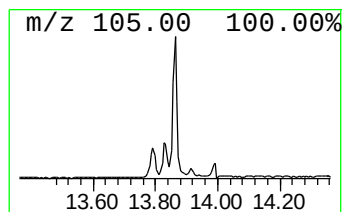
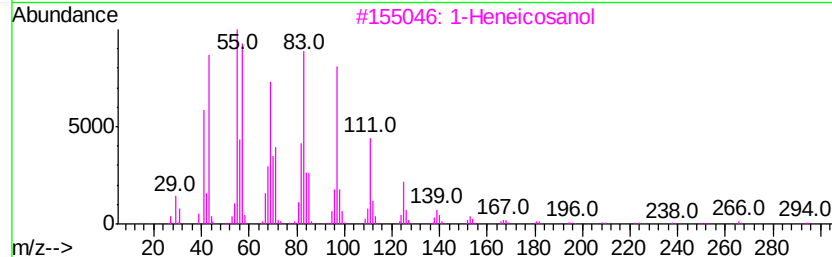
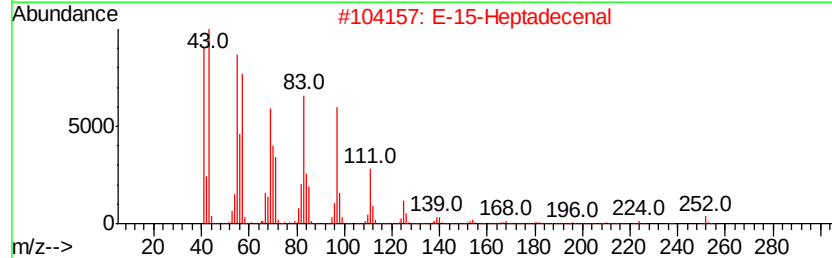
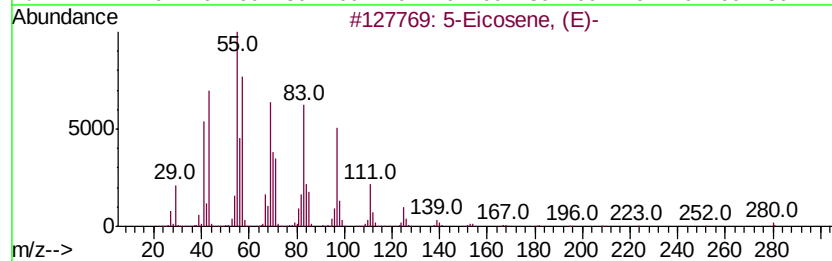
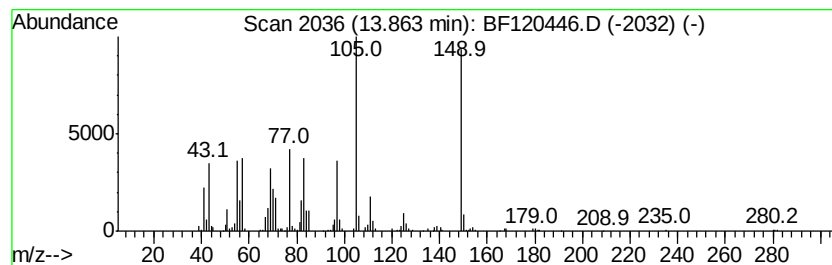
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 5 5-Eicosene, (E)- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.86	5.92 ng	516398	Chrysene-d12	13.99
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	5-Eicosene, (E)-	280 C20H40	074685-30-6	97
2	E-15-Heptadecenal	252 C17H32O	1000130-97-9	95
3	1-Heneicosanol	312 C21H44O	015594-90-8	93
4	Trifluoroacetoxy hexadecane	338 C18H33F3O2	006222-03-3	91
5	n-Nonadecanol-1	284 C19H40O	001454-84-8	89



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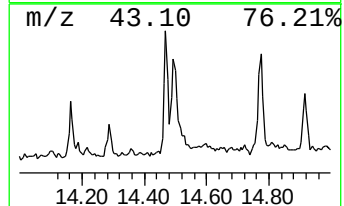
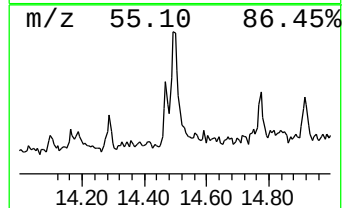
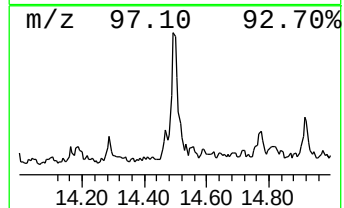
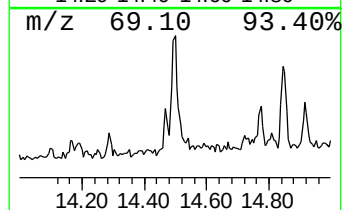
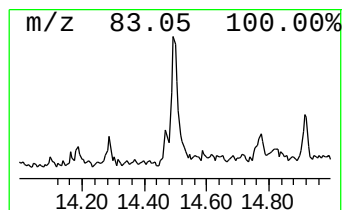
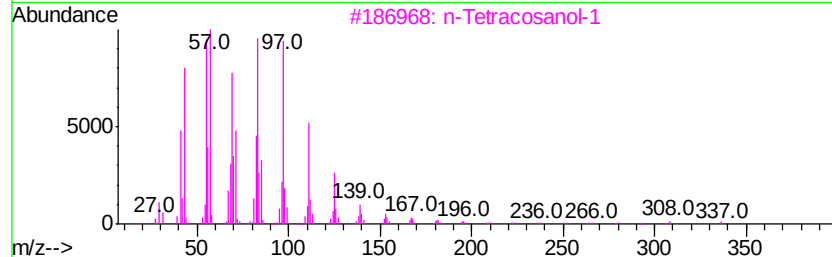
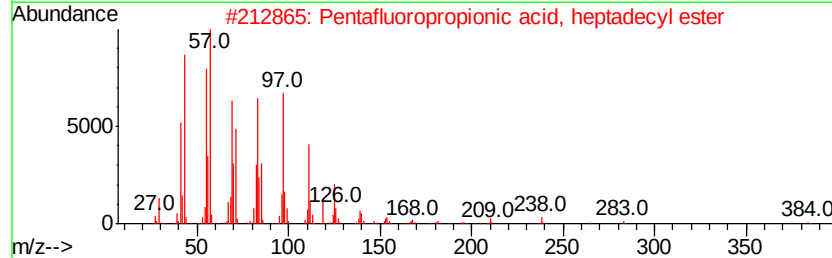
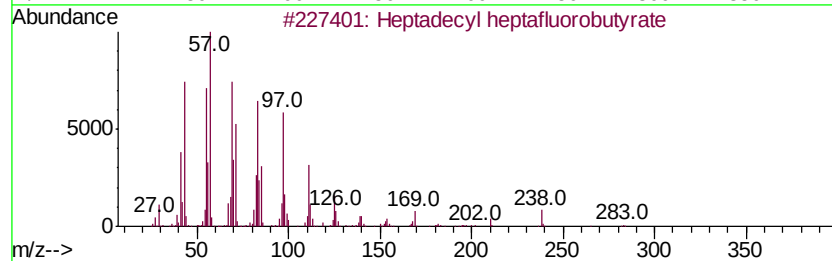
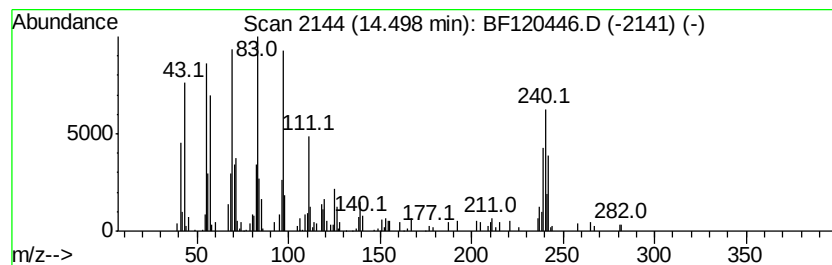
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 ClientSampleId :
 NM73-COMP-2020-0-6

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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 6 Heptadecyl heptafluorobutyrate Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.50	2.20 ng	192081	Chrysene-d12	13.99	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	62
2	Pentafluoropropionic acid, hepta...	402	C20H35F5O2	959218-78-1	50
3	n-Tetracosanol-1	354	C24H50O	000506-51-4	50
4	n-Heptadecanol-1	256	C17H36O	001454-85-9	45
5	1-Heptadecene	238	C17H34	006765-39-5	43



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF052920\
Data File : BF120446.D
Acq On : 29 May 2020 18:28
Operator : MA/SJ
Sample : L2773-20
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
NM73-COMP-2020-0-6

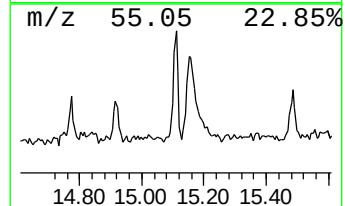
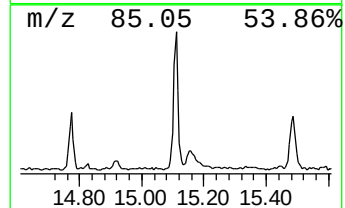
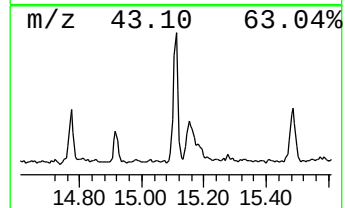
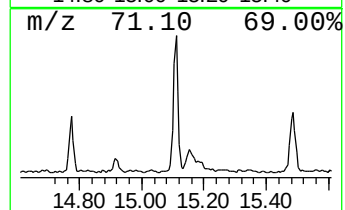
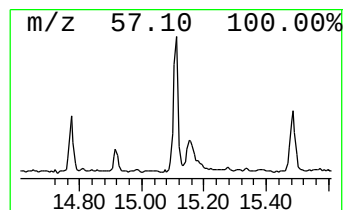
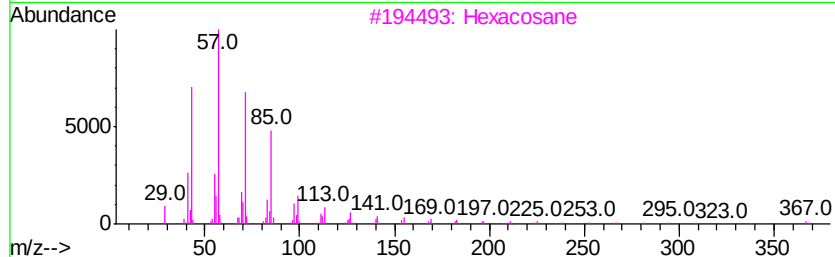
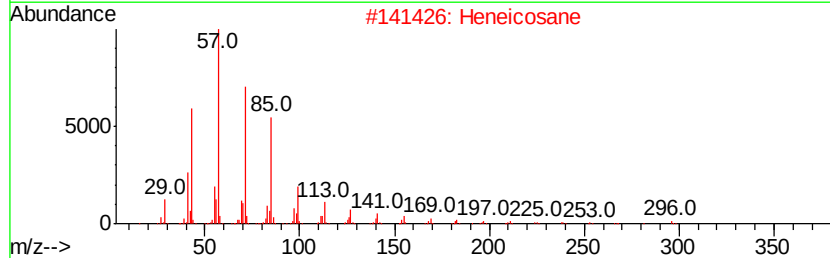
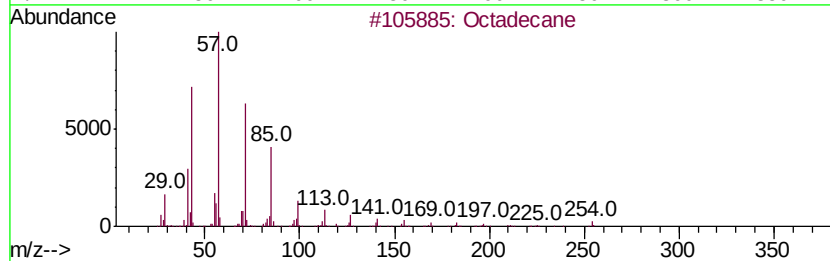
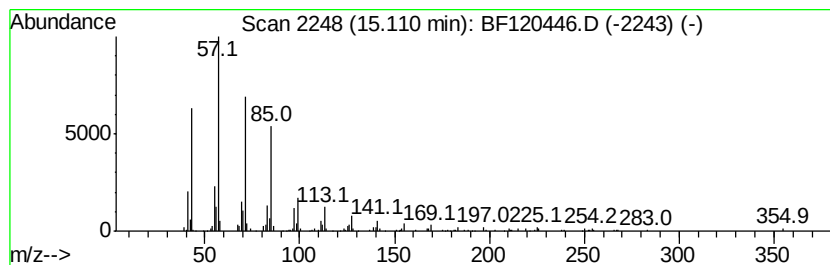
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Octadecane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.11	3.33 ng	256749	Perylene-d12	15.44

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecane	254	C18H38	000593-45-3	95
2		Heneicosane	296	C21H44	000629-94-7	90
3		Hexacosane	366	C26H54	000630-01-3	90
4		Heptadecane, 8-methyl-	254	C18H38	013287-23-5	90
5		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF052920\
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Misc :
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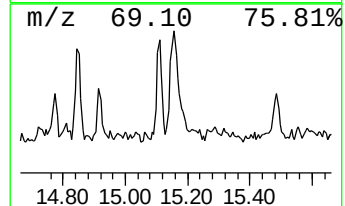
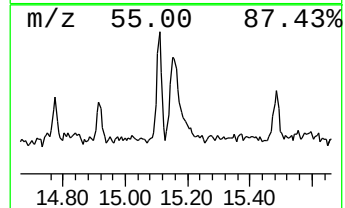
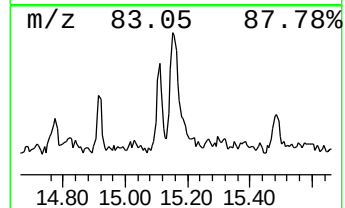
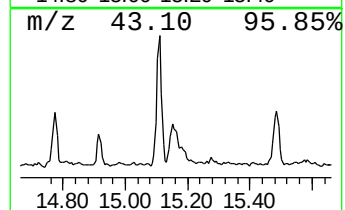
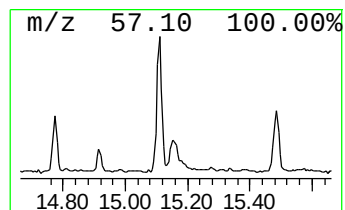
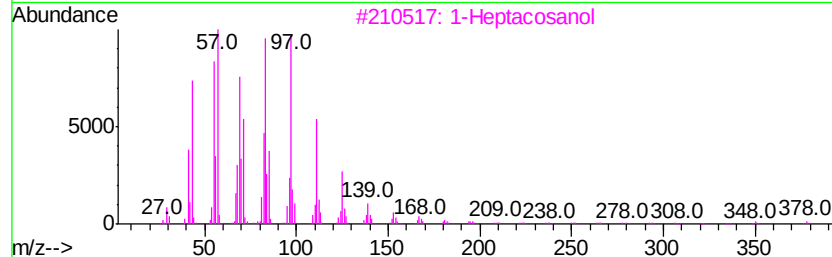
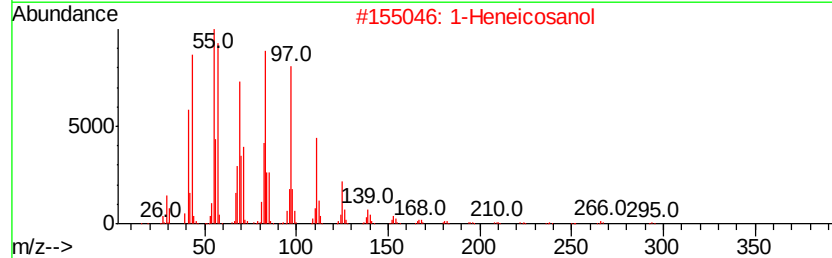
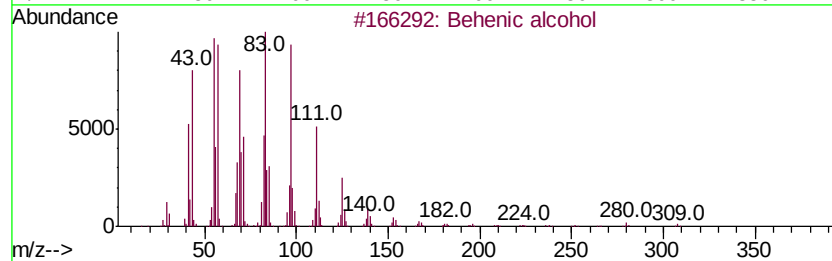
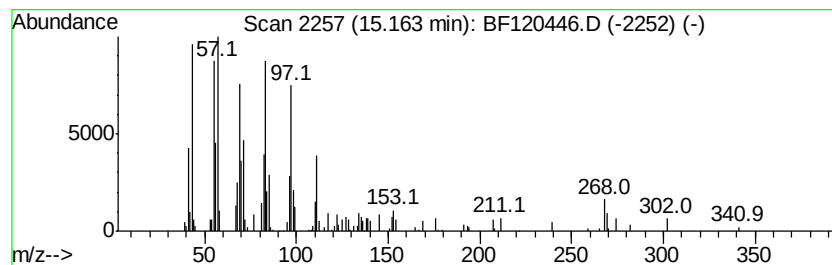
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Behenic alcohol Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.16	2.13 ng	164035	Perylene-d12	15.44
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Behenic alcohol	326 C22H46O	000661-19-8 93
2		1-Heneicosanol	312 C21H44O	015594-90-8 87
3		1-Heptacosanol	396 C27H56O	002004-39-9 87
4		Pentafluoropropionic acid, penta...	374 C18H31F5O2	959092-08-1 86
5		Heptafluorobutyric acid, pentade...	424 C19H31F7O2	959261-23-5 86



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF052920\
Data File : BF120446.D
Acq On : 29 May 2020 18:28
Operator : MA/SJ
Sample : L2773-20
Misc :
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Instrument :
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ClientSampleId :
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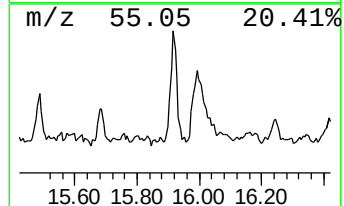
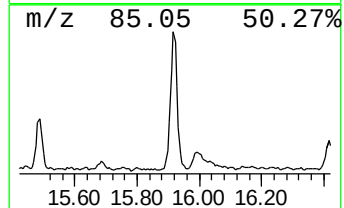
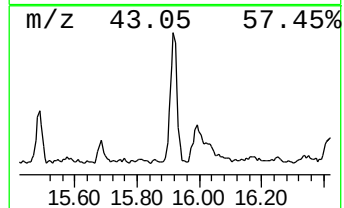
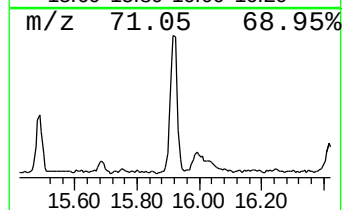
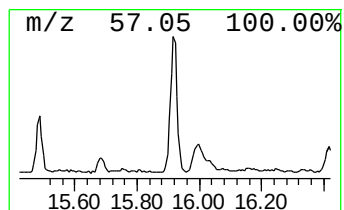
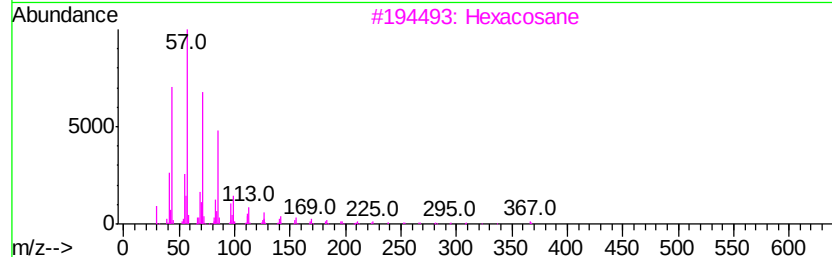
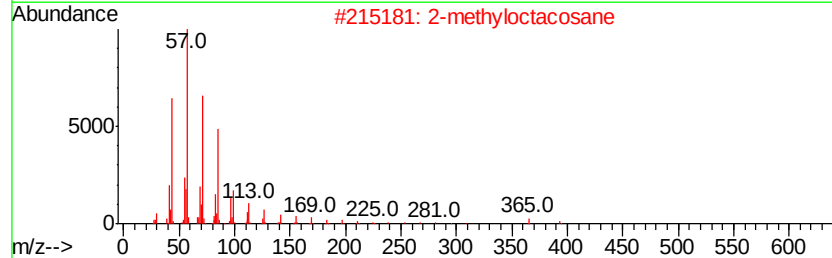
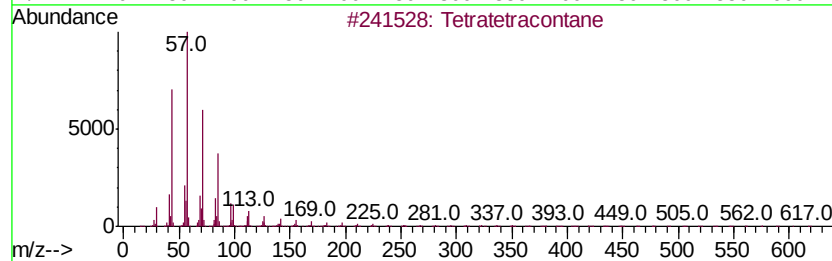
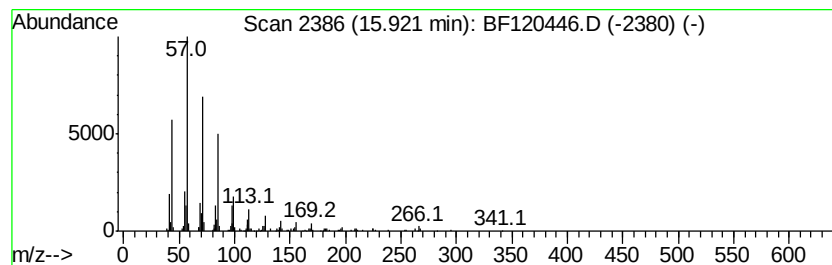
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Tetratetracontane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.92	4.11 ng	317121	Perylene-d12	15.44

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Tetratetracontane	619	C44H90	007098-22-8	90
2		2-methyloctacosane	408	C29H60	1000376-72-8	87
3		Hexacosane	366	C26H54	000630-01-3	87
4		Tetracosane	338	C24H50	000646-31-1	87
5		Hexadecane, 5-butyl-	282	C20H42	006912-07-8	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF052920\
Data File : BF120446.D
Acq On : 29 May 2020 18:28
Operator : MA/SJ
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Misc :
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ClientSampleId :
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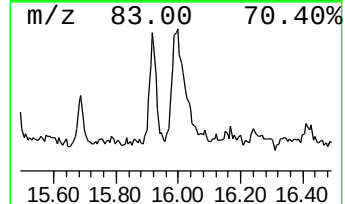
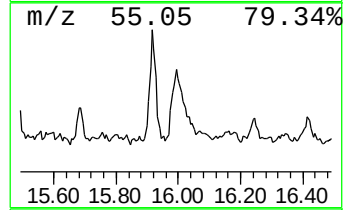
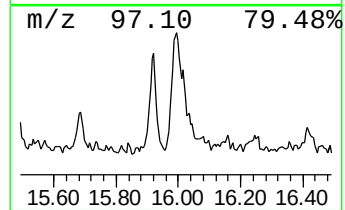
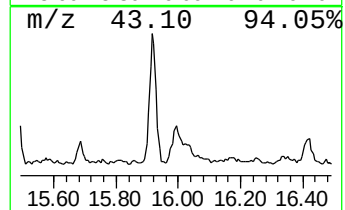
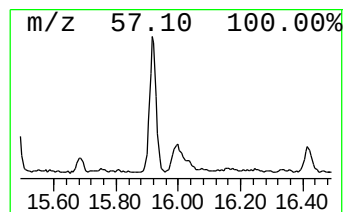
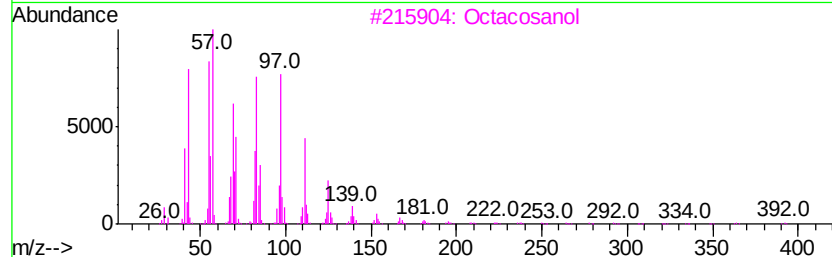
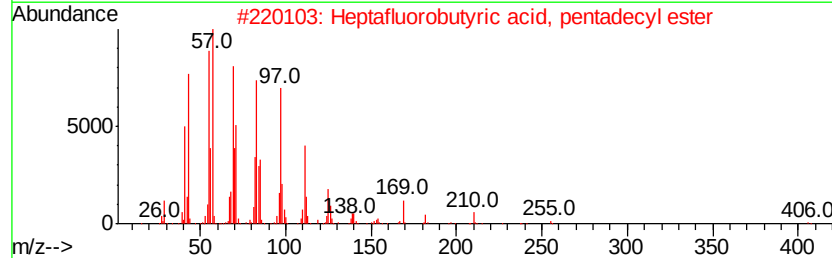
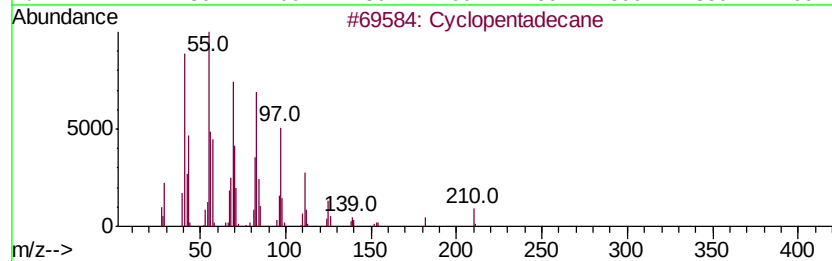
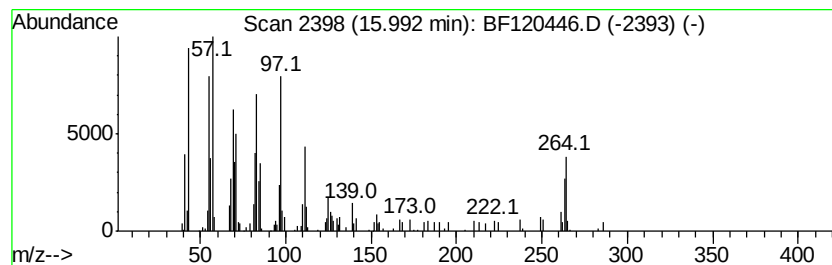
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Cyclopentadecane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.99	2.83 ng	218642	Perylene-d12	15.44

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopentadecane	210	C15H30	000295-48-7	93
2		Heptafluorobutyric acid, pentade...	424	C19H31F7O2	959261-23-5	89
3		Octacosanol	410	C28H58O	000557-61-9	89
4		Pentafluoropropionic acid, penta...	374	C18H31F5O2	959092-08-1	89
5		Dichloroacetic acid, heptadecyl ...	366	C19H36Cl2O2	1000282-98-2	86



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF052920\
Data File : BF120446.D
Acq On : 29 May 2020 18:28
Operator : MA/SJ
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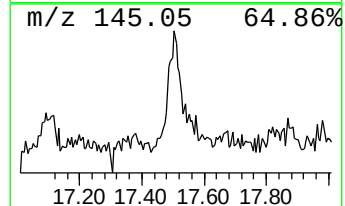
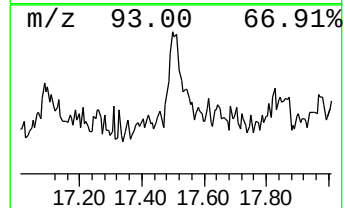
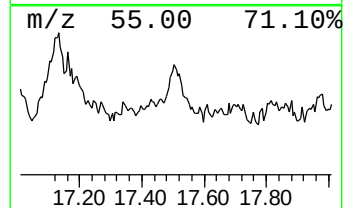
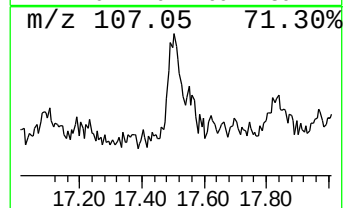
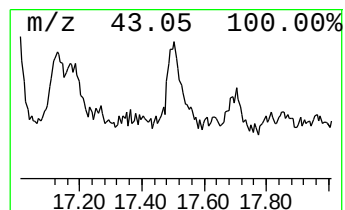
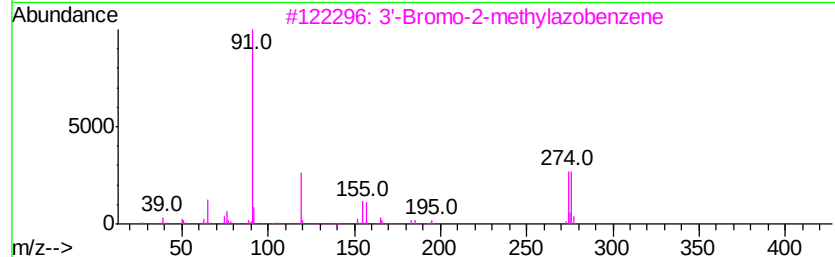
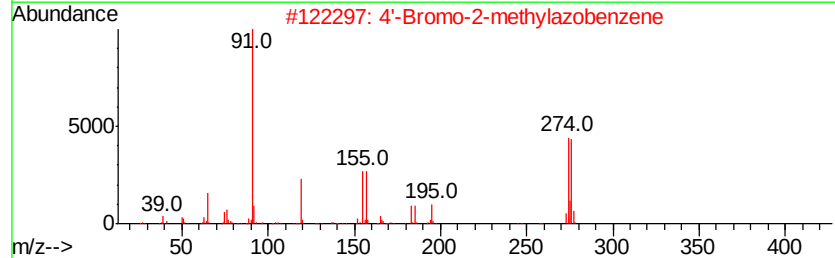
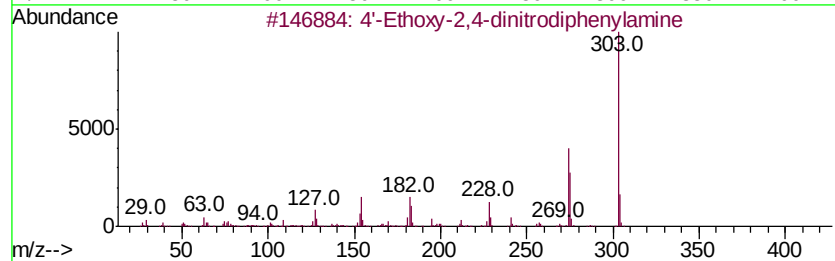
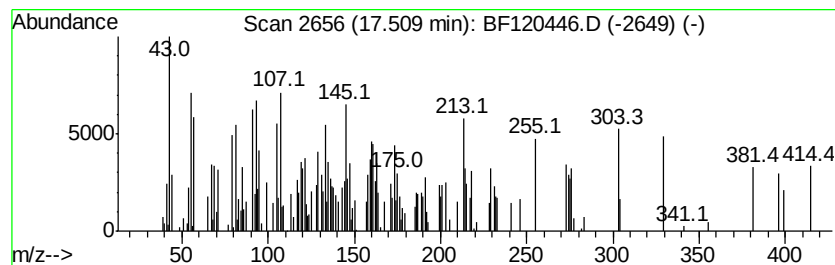
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 unknown17.51 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.51	3.76 ng	290072	Perylene-d12	15.44

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4'-Ethoxy-2,4-dinitrodiphenylamine	303	C14H13N3O5	006943-24-4	12
2		4'-Bromo-2-methylazobenzene	274	C13H11BrN2	092022-22-5	11
3		3'-Bromo-2-methylazobenzene	274	C13H11BrN2	092022-18-9	11
4		N-(4-Bromobenzylidene)-p-toluidine	273	C14H12BrN	039128-27-3	11
5		1H-Pyrazolo[3,4-b]quinoline, 6-e...	303	C19H17N3O	1000304-74-4	10



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF052920\
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 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
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Methylene chloride	1.96	32.9	ng	2623020	1	6.83	1594270	20.0
Butane, 2-methoxy...	2.11	18.8	ng	1499540	1	6.83	1594270	20.0
5-Eicosene, (E)-	13.86	5.9	ng	516398	5	13.99	1743850	20.0
Heptadecyl hepta...	14.50	2.2	ng	192081	5	13.99	1743850	20.0
Octadecane	15.11	3.3	ng	256749	6	15.44	1542460	20.0
Behenic alcohol	15.16	2.1	ng	164035	6	15.44	1542460	20.0
Tetratetracontane	15.92	4.1	ng	317121	6	15.44	1542460	20.0
Cyclopentadecane	15.99	2.8	ng	218642	6	15.44	1542460	20.0
unknown17.51	17.51	3.8	ng	290072	6	15.44	1542460	20.0