

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP011720\
 Data File : BP001658.D
 Acq On : 17 Jan 2020 15:39
 Operator : JU
 Sample : L1171-01
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BW-1A-20.0

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_P\METHODS\8270-BP010820.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	2.864	9	13	22	rBV2	31765	62611	2.81%	0.368%
2	2.946	22	27	36	rVB	227153	283005	12.70%	1.665%
3	4.793	337	341	350	rVB	179898	252254	11.32%	1.484%
4	5.264	415	421	430	rBV	782745	1130504	50.72%	6.653%
5	6.834	681	688	704	rBV	875026	1419359	63.68%	8.353%
6	7.175	739	746	756	rBV	839783	1293914	58.06%	7.615%
7	7.628	817	823	833	rVB	202065	317541	14.25%	1.869%
8	7.952	871	878	886	rBV	595337	940400	42.19%	5.534%
9	8.781	1012	1019	1036	rBV	524936	868729	38.98%	5.112%
10	10.404	1287	1295	1307	rBV	228599	406719	18.25%	2.393%
11	12.898	1711	1719	1739	rBV	1123038	1659451	74.46%	9.766%
12	13.745	1857	1863	1876	rBV	66469	108656	4.88%	0.639%
13	14.281	1947	1954	1967	rBV2	392785	585878	26.29%	3.448%
14	15.781	2203	2209	2236	rBV	916300	1383598	62.08%	8.142%
15	17.039	2417	2423	2438	rBV	460176	705275	31.64%	4.150%
16	19.628	2857	2863	2870	rBV	1880296	2228768	100.00%	13.116%
17	19.816	2889	2895	2910	rVB3	30580	142071	6.37%	0.836%
18	20.839	3065	3069	3072	rBV	15996	23570	1.06%	0.139%
19	20.886	3072	3077	3083	rVV2	186402	288082	12.93%	1.695%
20	20.939	3083	3086	3094	rVB	33201	51604	2.32%	0.304%
21	21.145	3116	3121	3130	rBV	430384	567903	25.48%	3.342%
22	21.322	3147	3151	3163	rVB2	116118	153054	6.87%	0.901%
23	21.757	3221	3225	3230	rBV	183235	219897	9.87%	1.294%
24	22.221	3300	3304	3317	rVB	272805	415932	18.66%	2.448%
25	22.739	3386	3392	3402	rVB	250199	381946	17.14%	2.248%
26	23.310	3484	3489	3494	rBV	198899	320833	14.40%	1.888%
27	23.374	3494	3500	3507	rVB	211553	402712	18.07%	2.370%
28	23.968	3595	3601	3607	rVB	124334	244916	10.99%	1.441%
29	24.727	3724	3730	3737	rVB2	64167	133536	5.99%	0.786%

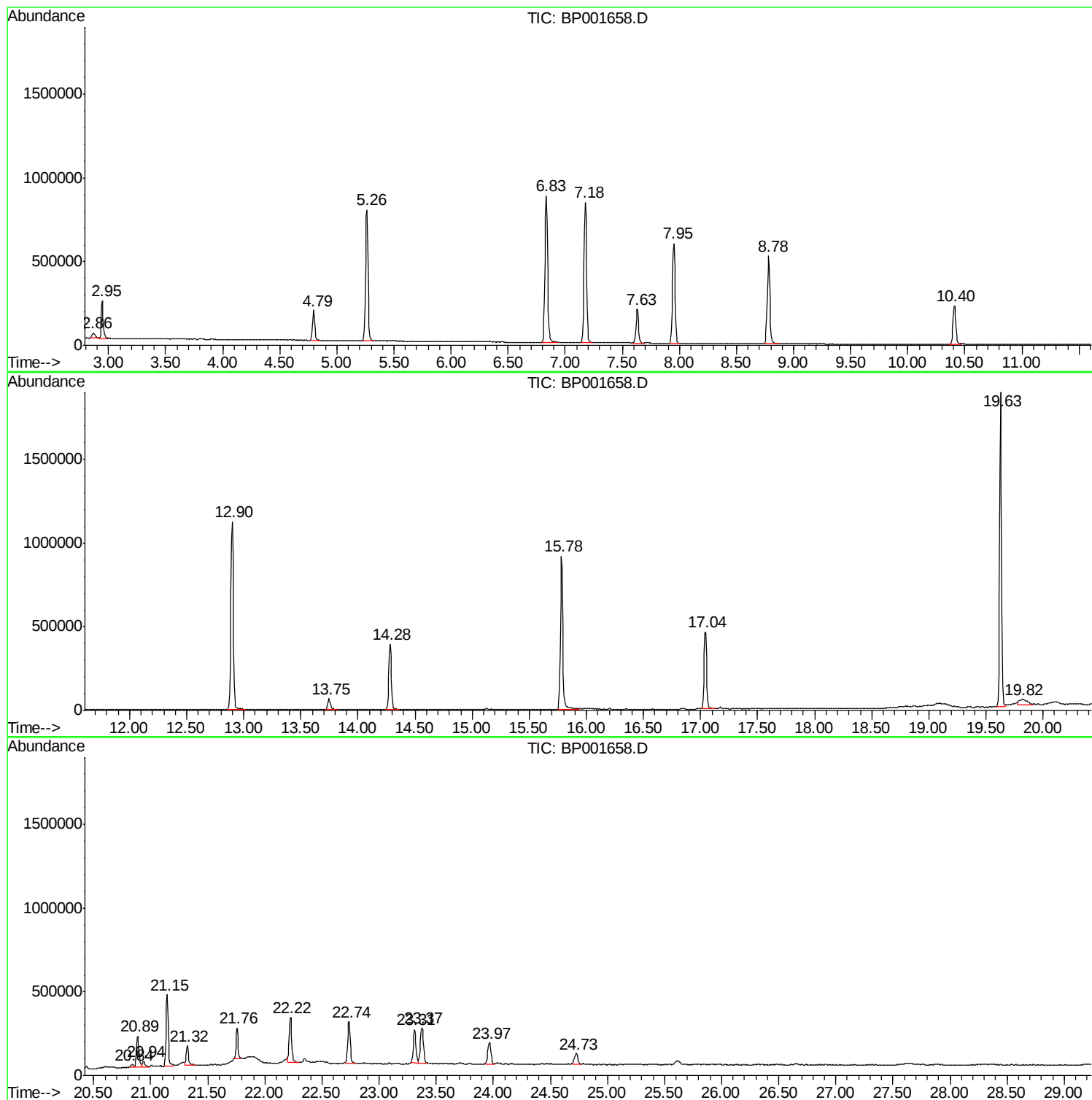
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP010820.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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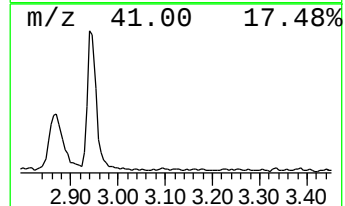
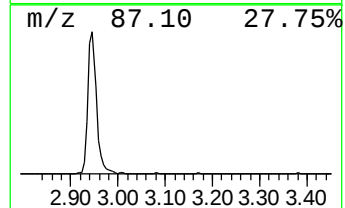
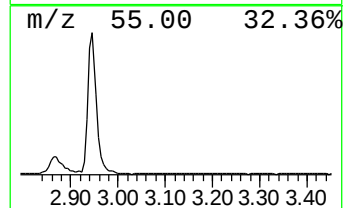
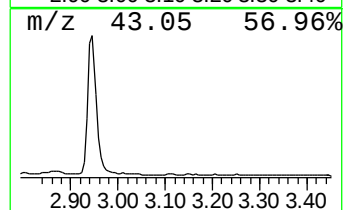
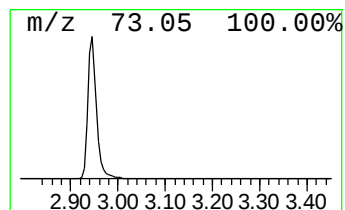
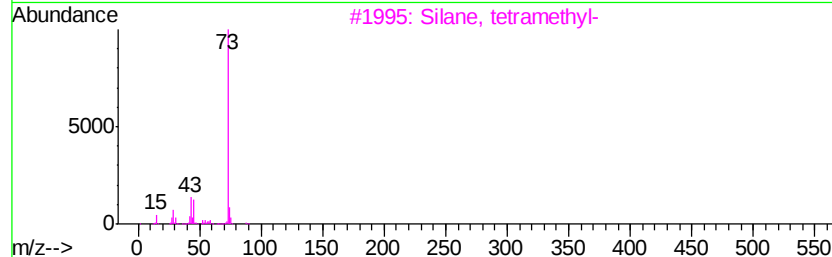
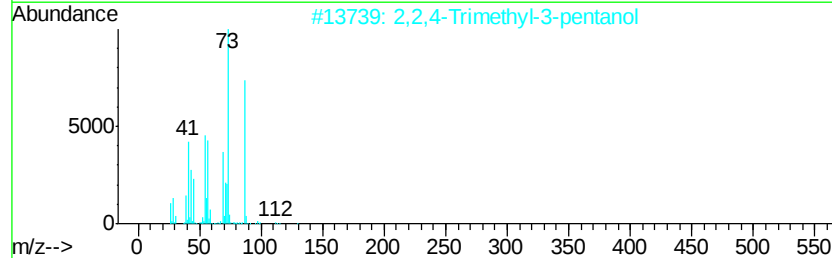
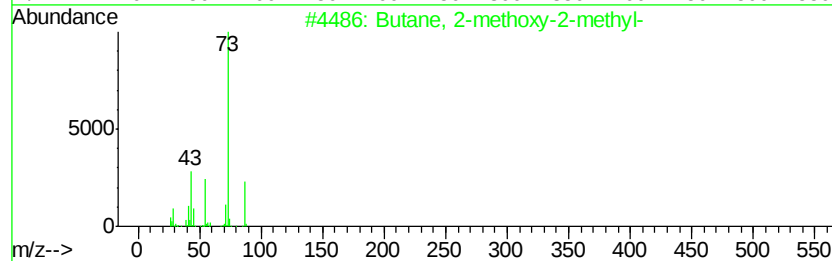
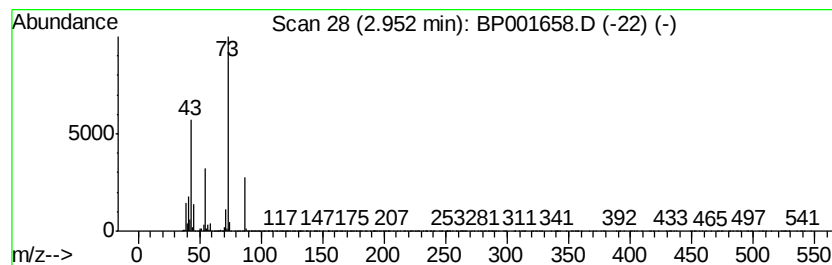
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 Butane, 2-methoxy-2-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.95	17.82 ng	283005	1,4-Dichlorobenzene-d4	7.63

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	40
3		Silane, tetramethyl-	88	C4H12Si	000075-76-3	32
4		1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	9
5		1,4-Butanediamine	88	C4H12N2	000110-60-1	9



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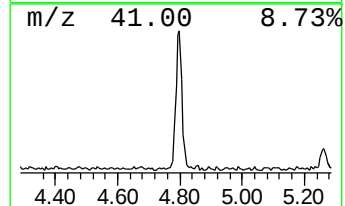
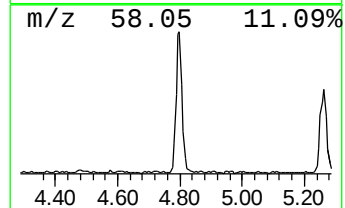
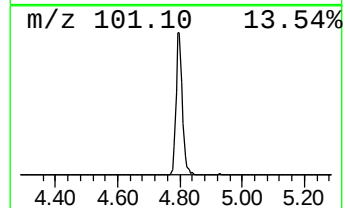
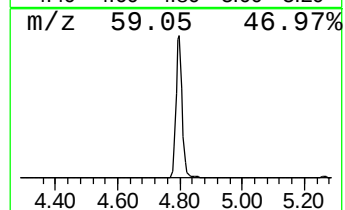
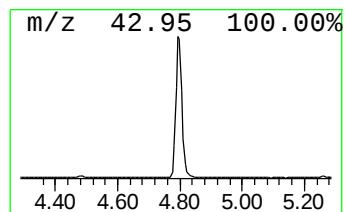
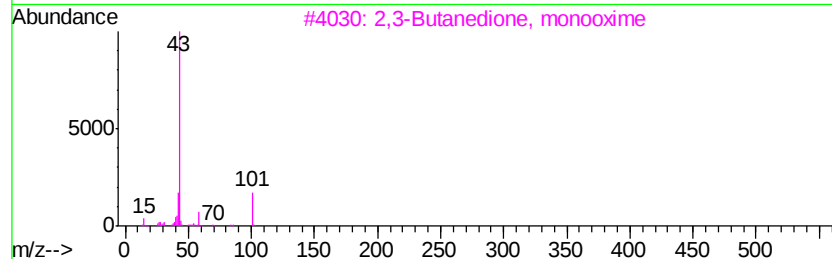
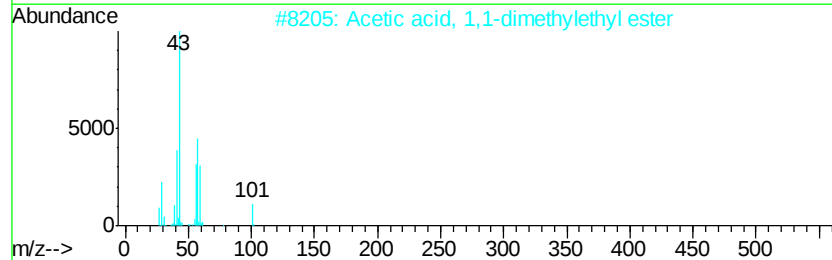
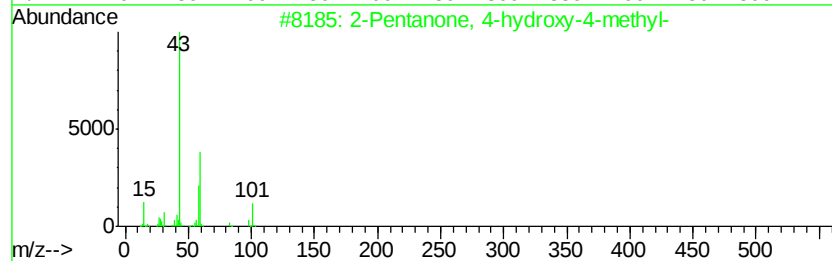
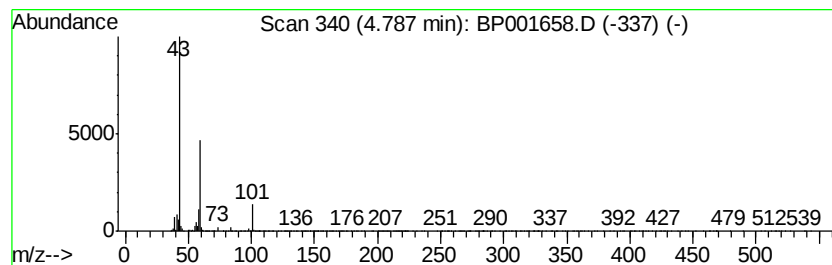
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TIC Integration Parameters: LSCINT.P

Peak Number 3 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.79	15.89 ng	252254	1,4-Dichlorobenzene-d4	7.63

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	72
2		Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	28
3		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	27
4		1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10
5		2-Propanone, 1-(1-methylethoxy)-	116	C6H12O2	042781-12-4	9



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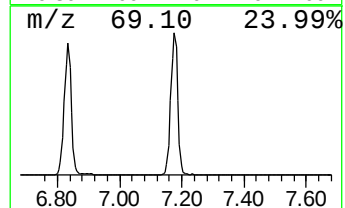
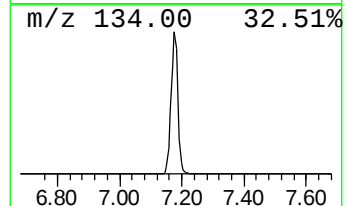
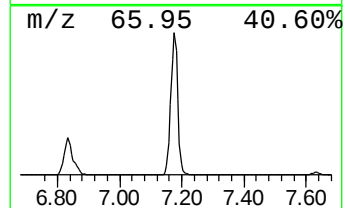
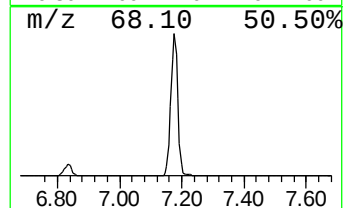
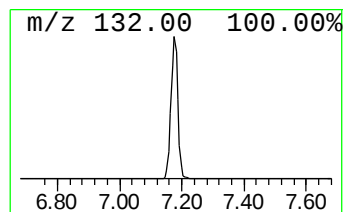
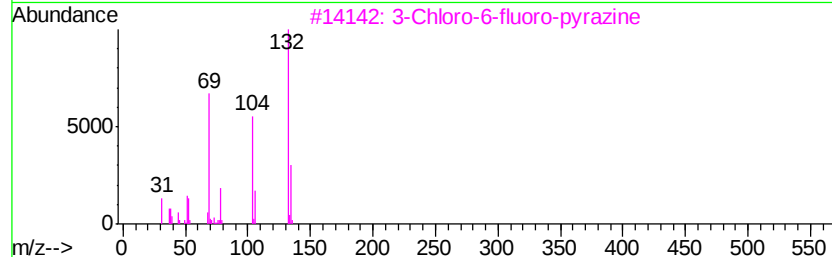
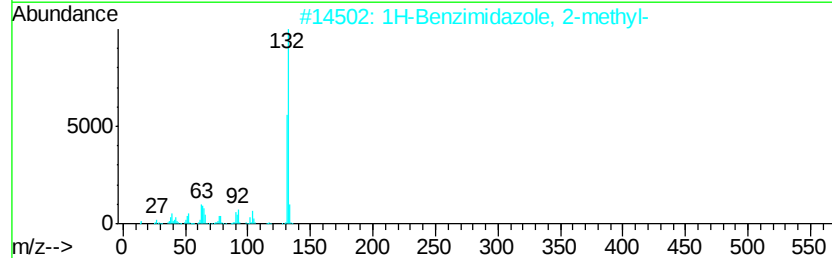
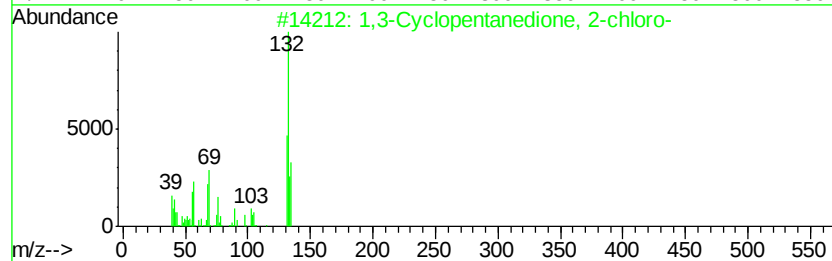
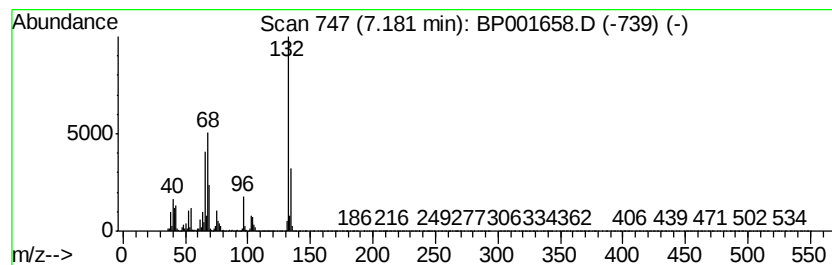
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Peak Number 4 unknown7.18 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.18	81.50 ng	1293910	1,4-Dichlorobenzene-d4	7.63

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,3-Cyclopentanedione, 2-chloro-	132	C5H5ClO2	014203-19-1	22
2		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	14
3		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	14
4		Pyrazolo(2,3-a)pyridine, 7-methyl-	132	C8H8N2	034760-58-2	14
5		1,2,4-Trifluorobenzene	132	C6H3F3	000367-23-7	10



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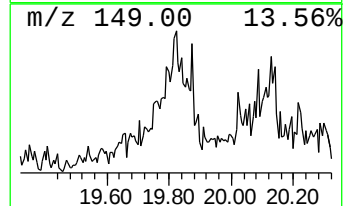
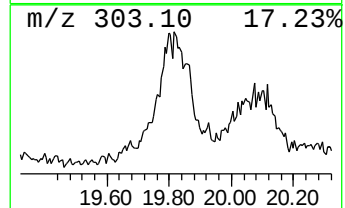
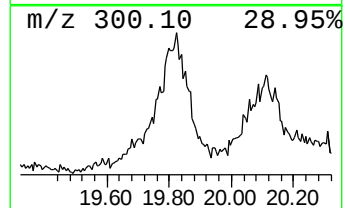
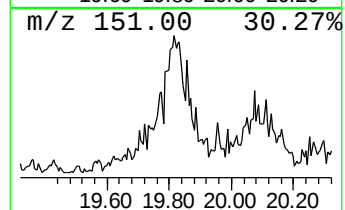
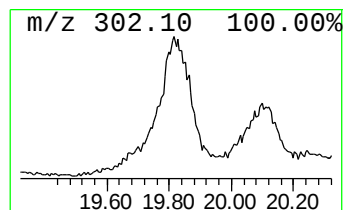
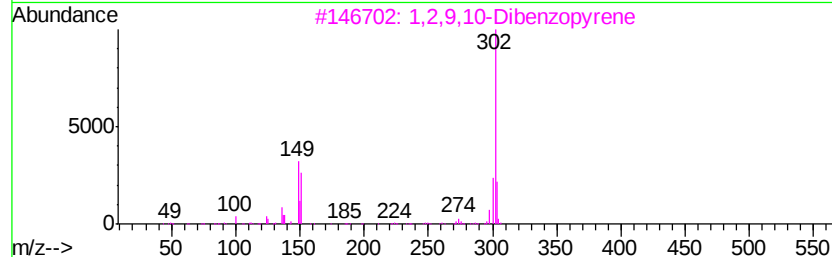
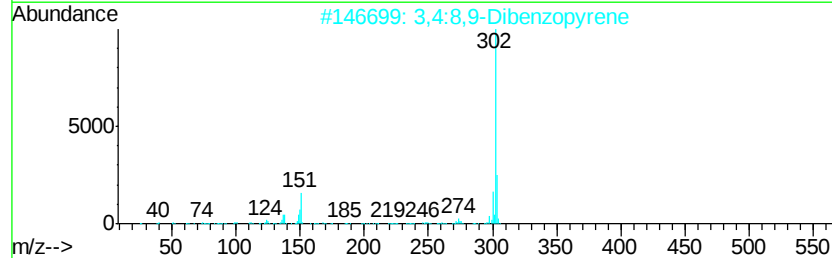
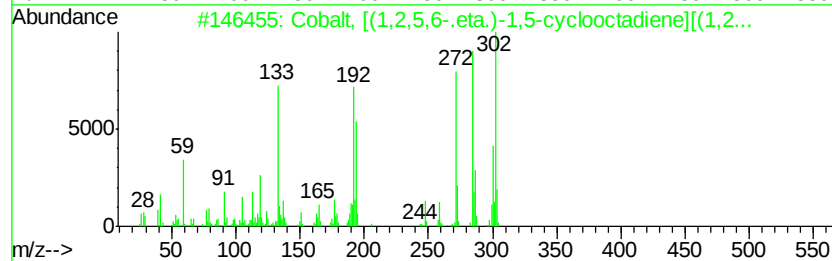
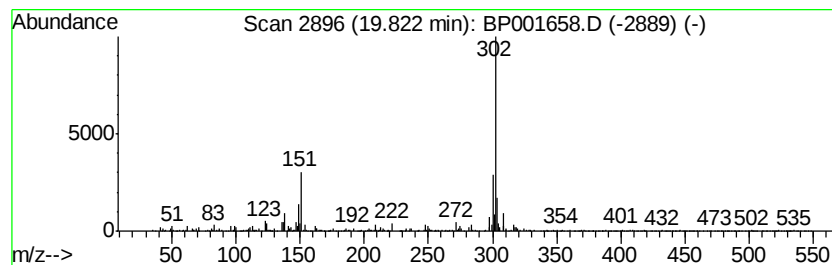
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Peak Number 6 Cobalt, [(1,2,5,6-.eta.)-1,... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.82	5.00 ng	142071	Chrysene-d12	21.15

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cobalt, [(1,2,5,6-.eta.)-1,5-cvc...	302	C18H27Co	082595-75-3	62
2		3,4:8,9-Dibenzopyrene	302	C24H14	000189-64-0	58
3		1,2,9,10-Dibenzopyrene	302	C24H14	000191-30-0	58
4		10,11-Dihydro-10-hydroxy-2,3,6-t...	302	C17H18O5	019172-35-1	56
5		Coronene, 1,2-dihydro-	302	C24H14	107716-56-3	53



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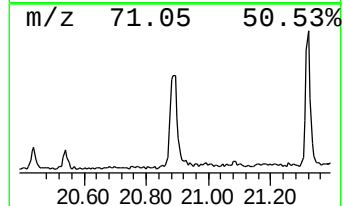
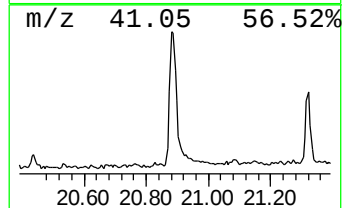
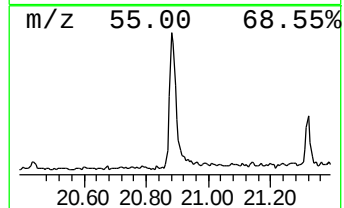
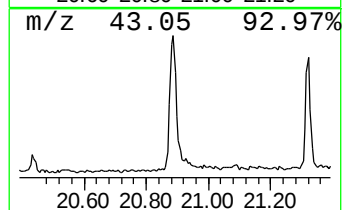
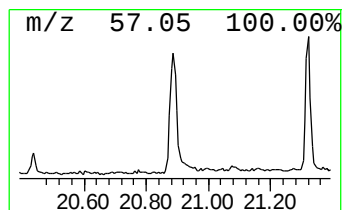
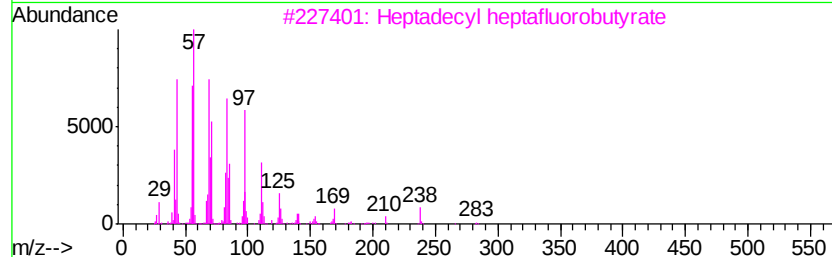
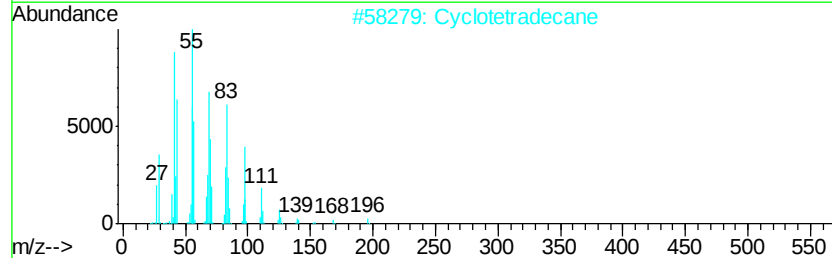
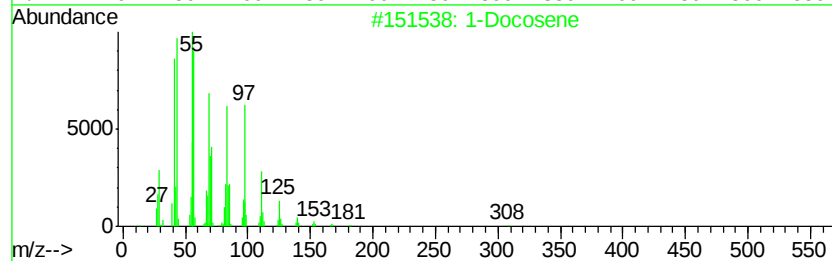
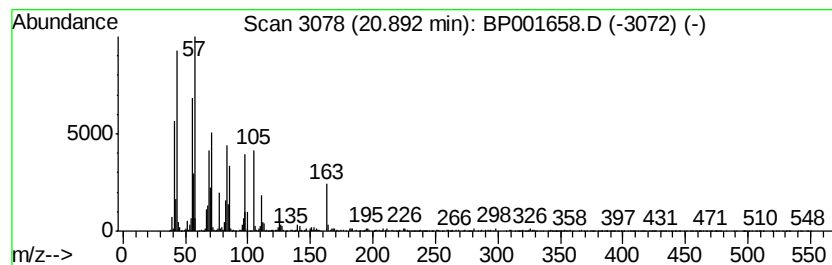
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Peak Number 7 1-Docosene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.89	10.15 ng	288082	Chrysene-d12	21.15

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Docosene		308	C22H44	001599-67-3	93
2	Cyclotetradecane		196	C14H28	000295-17-0	91
3	Heptadecyl heptafluorobutvrate		452	C21H35F7O2	959085-66-6	90
4	5-Eicosene, (E)-		280	C20H40	074685-30-6	89
5	Cyclohexadecane, 1,2-diethyl-		280	C20H40	1000155-85-3	87



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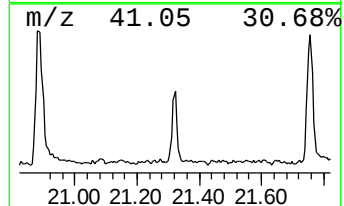
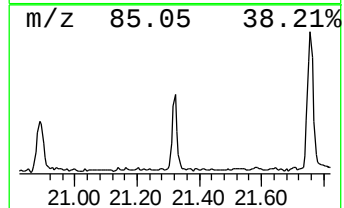
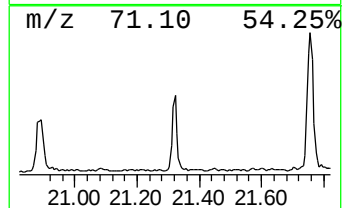
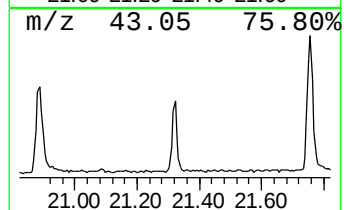
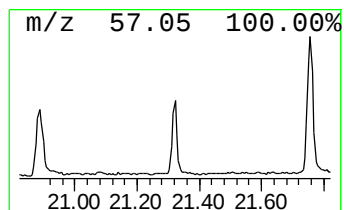
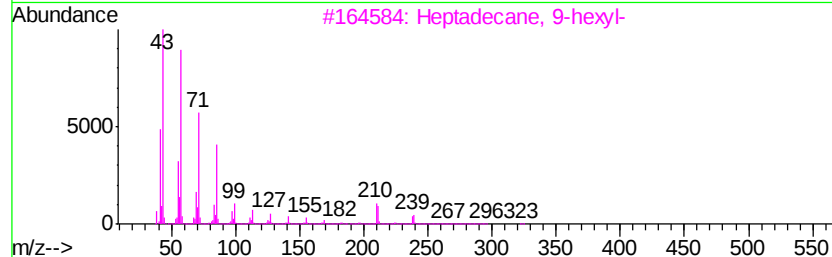
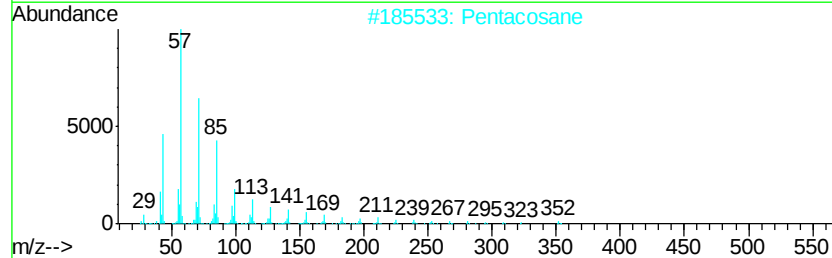
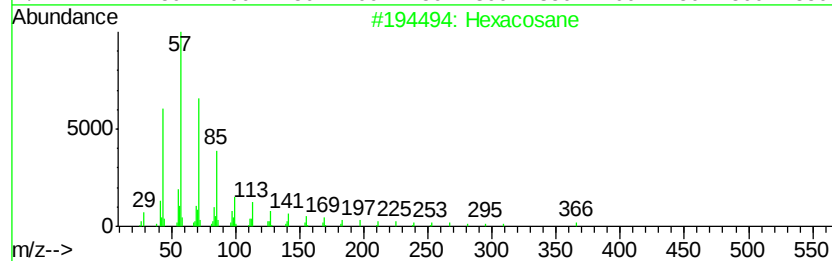
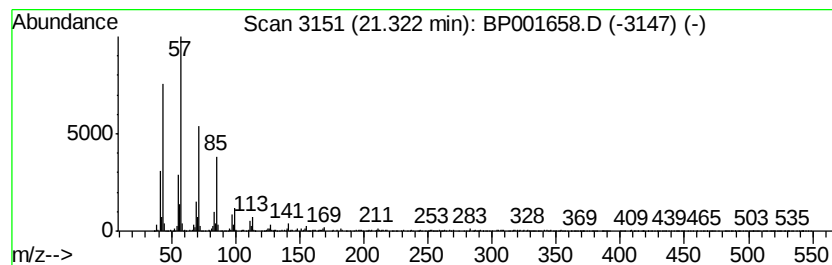
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Hexacosane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.32	5.39 ng	153054	Chrysene-d12	21.15

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexacosane	366	C26H54	000630-01-3	94
2		Pentacosane	352	C25H52	000629-99-2	93
3		Heptadecane, 9-hexyl-	324	C23H48	055124-79-3	91
4		Heptacosane	380	C27H56	000593-49-7	91
5		Octacosane	394	C28H58	000630-02-4	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP011720\
 Data File : BP001658.D
 Acq On : 17 Jan 2020 15:39
 Operator : JU
 Sample : L1171-01
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

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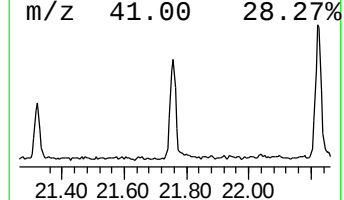
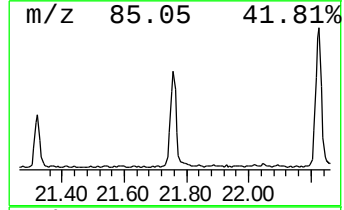
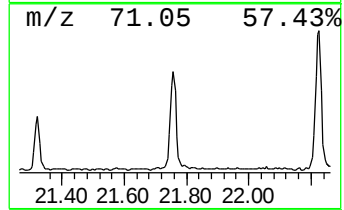
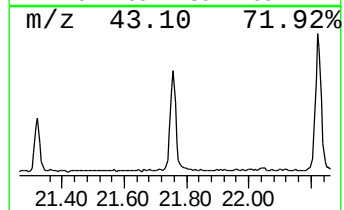
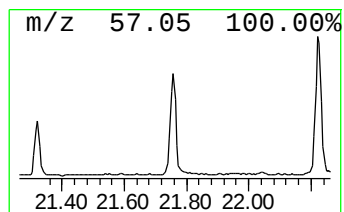
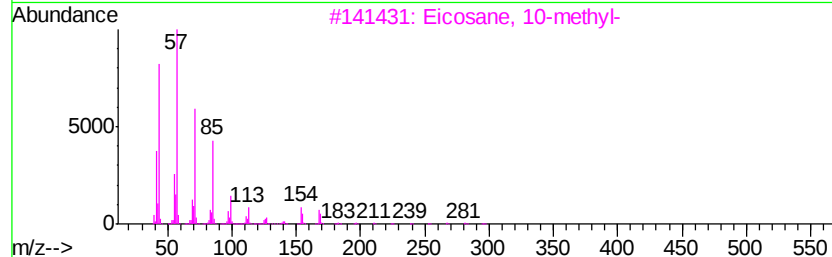
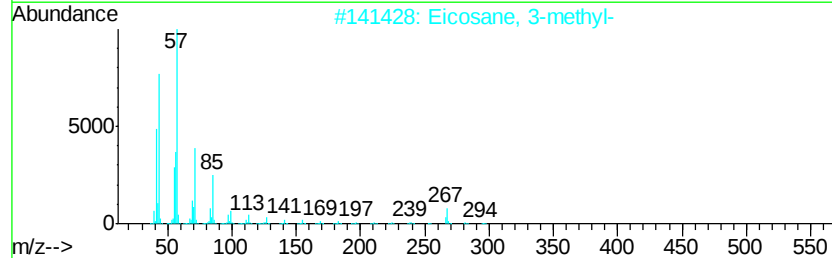
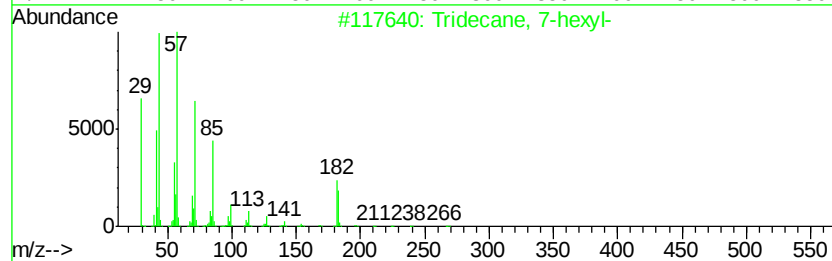
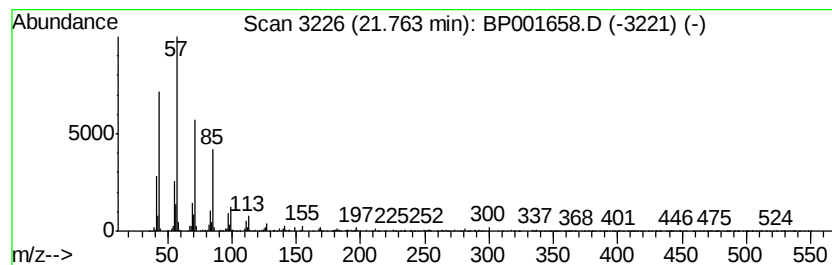
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

 Peak Number 9 Tridecane, 7-hexyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.76	7.74 ng	219897	Chrysene-d12	21.15

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tridecane, 7-hexyl-	268	C19H40	007225-66-3	96
2		Eicosane, 3-methyl-	296	C21H44	006418-46-8	93
3		Eicosane, 10-methyl-	296	C21H44	054833-23-7	90
4		Heptacosane	380	C27H56	000593-49-7	90
5		Heneicosane	296	C21H44	000629-94-7	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP011720\
Data File : BP001658.D
Acq On : 17 Jan 2020 15:39
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Sample : L1171-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

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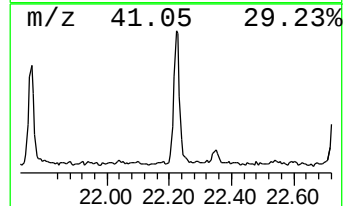
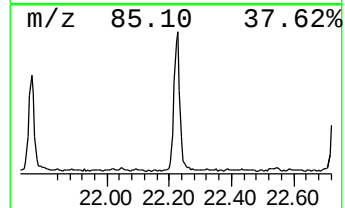
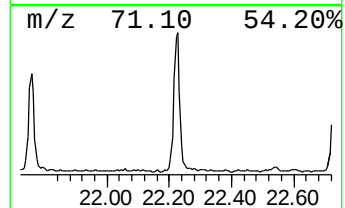
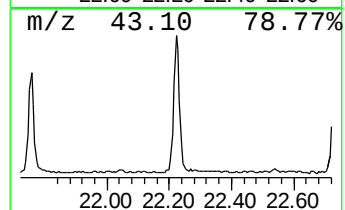
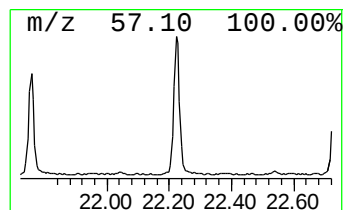
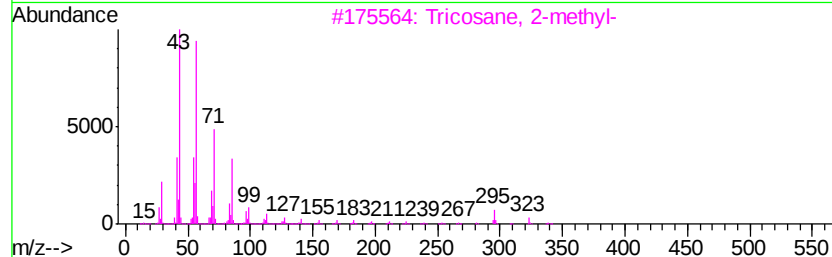
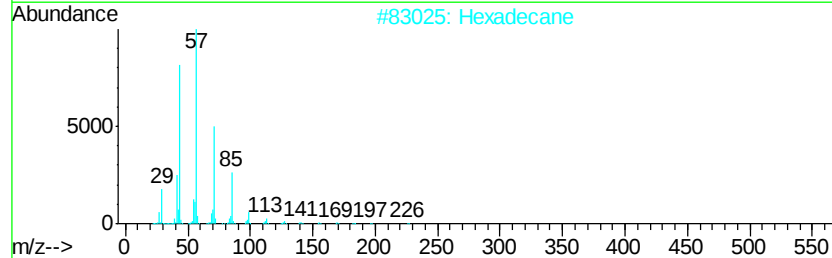
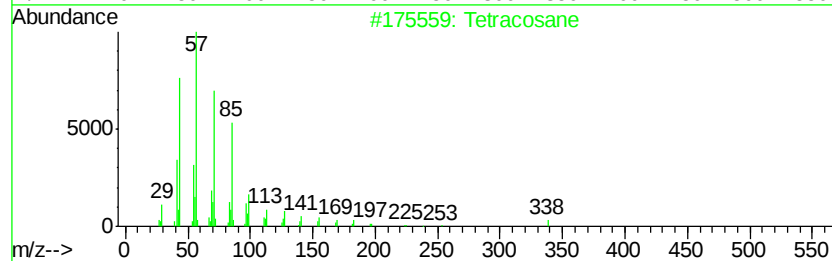
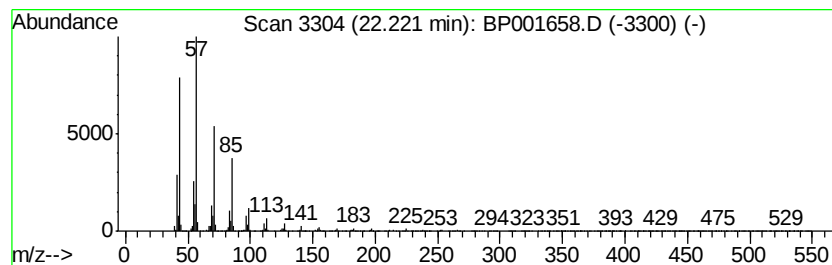
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Tetracosane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.22	14.65 ng	415932	Chrysene-d12	21.15

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetracosane	338	C24H50	000646-31-1	95
2		Hexadecane	226	C16H34	000544-76-3	94
3		Tricosane, 2-methyl-	338	C24H50	001928-30-9	91
4		Hexatriacontane	507	C36H74	000630-06-8	91
5		Octacosane	394	C28H58	000630-02-4	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP011720\
Data File : BP001658.D
Acq On : 17 Jan 2020 15:39
Operator : JU
Sample : L1171-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

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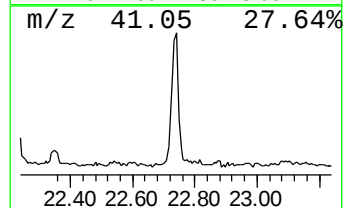
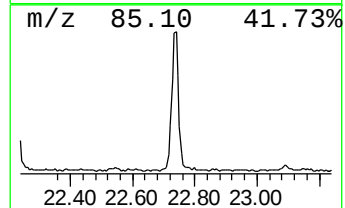
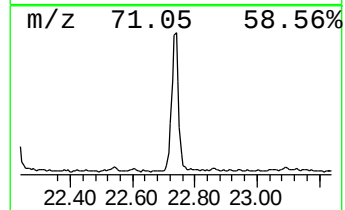
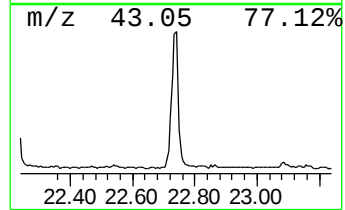
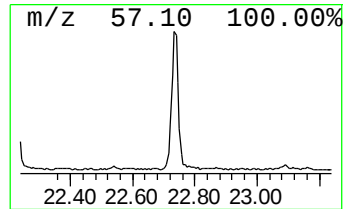
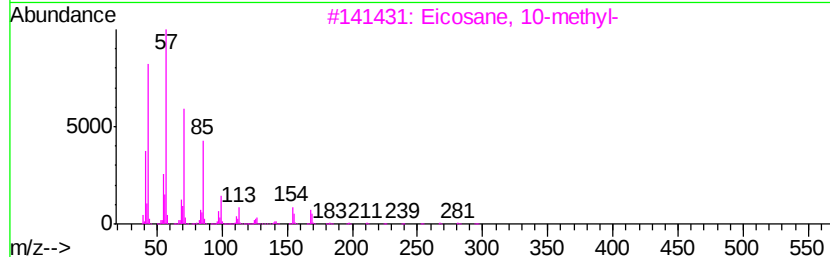
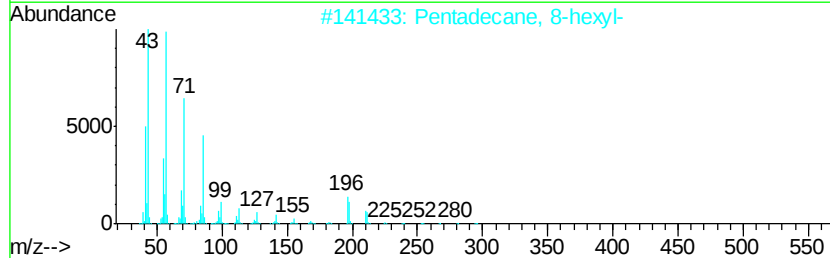
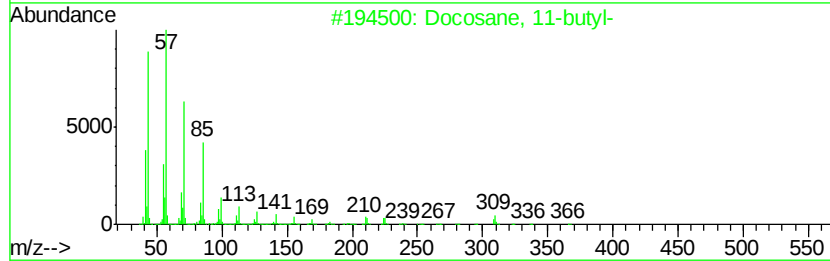
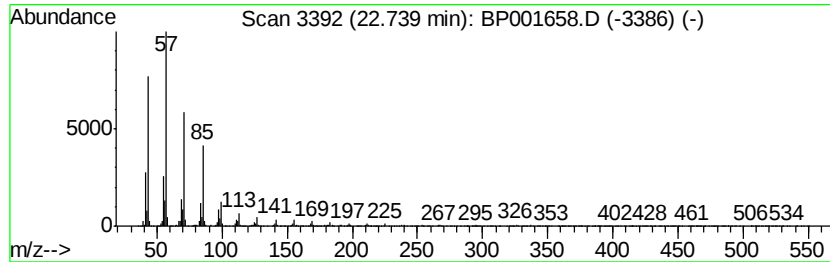
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Docosane, 11-butyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.74	18.97 ng	381946	Perylene-d12	23.37

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Docosane, 11-butyl-	366	C26H54	013475-76-8	95
2		Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	94
3		Eicosane, 10-methyl-	296	C21H44	054833-23-7	93
4		Tricosane	324	C23H48	000638-67-5	93
5		Heptacosane	380	C27H56	000593-49-7	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP011720\
Data File : BP001658.D
Acq On : 17 Jan 2020 15:39
Operator : JU
Sample : L1171-01
Misc :
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Instrument :
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ClientSampleId :
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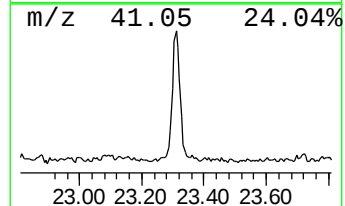
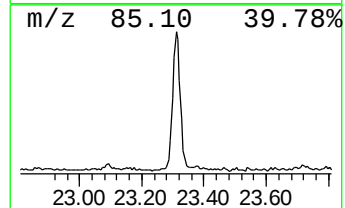
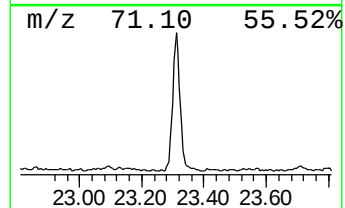
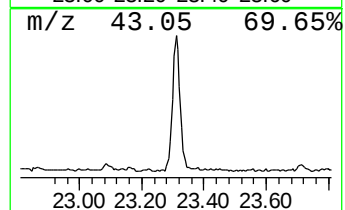
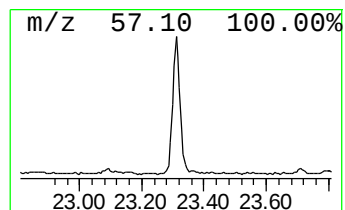
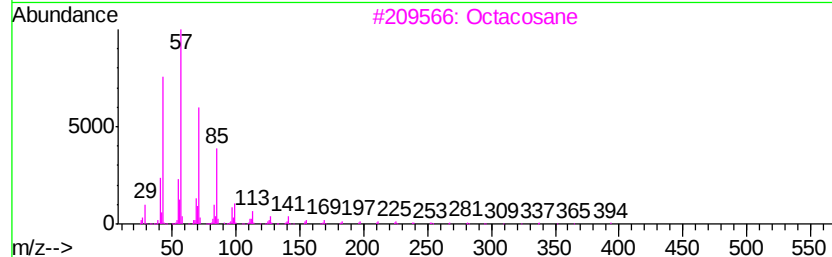
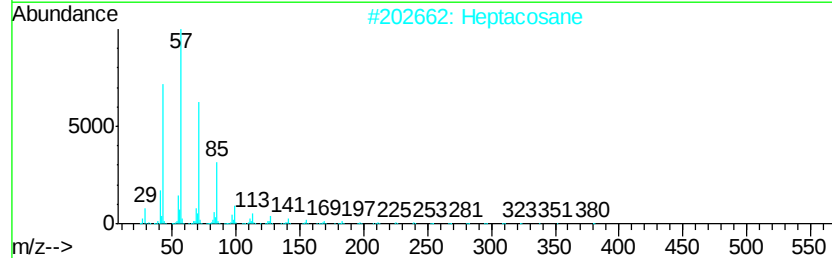
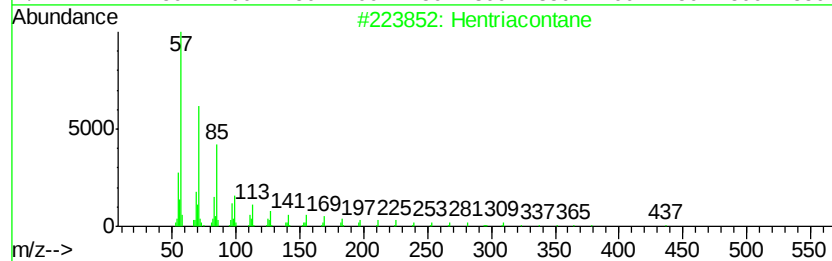
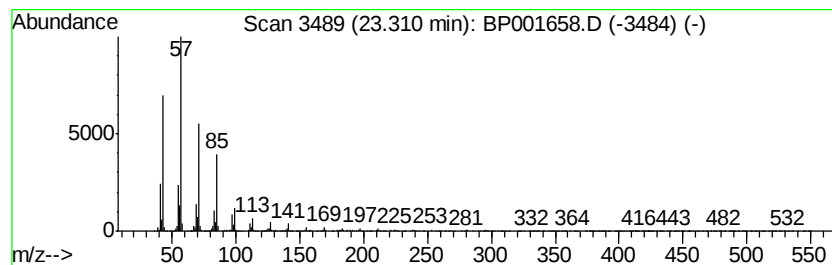
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 12 Hentriacontane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.31	15.93 ng	320833	Perylene-d12	23.37

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hentriacontane	437	C31H64	000630-04-6	91
2		Heptacosane	380	C27H56	000593-49-7	91
3		Octacosane	394	C28H58	000630-02-4	91
4		Tetracosane	338	C24H50	000646-31-1	91
5		Hexatriacontane	507	C36H74	000630-06-8	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP011720\
Data File : BP001658.D
Acq On : 17 Jan 2020 15:39
Operator : JU
Sample : L1171-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

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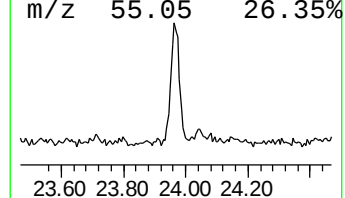
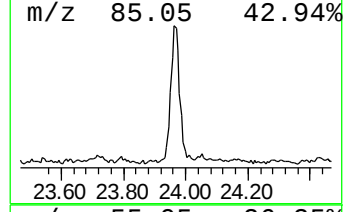
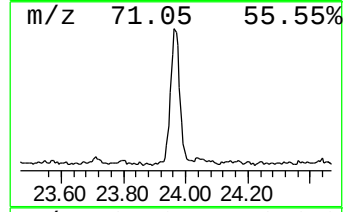
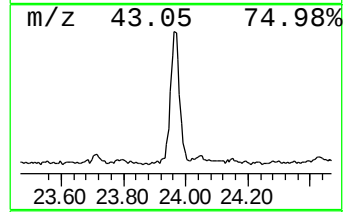
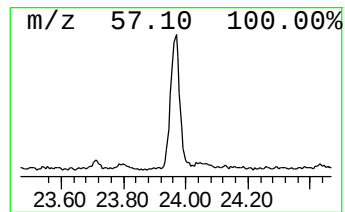
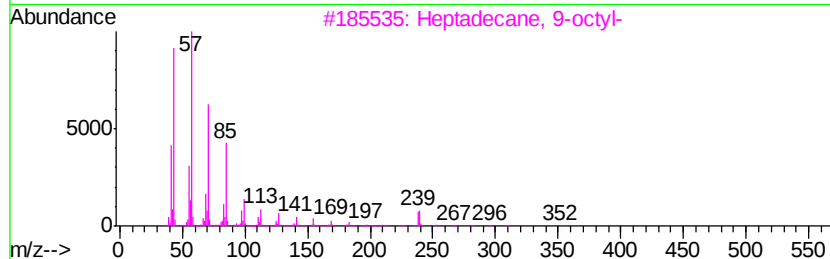
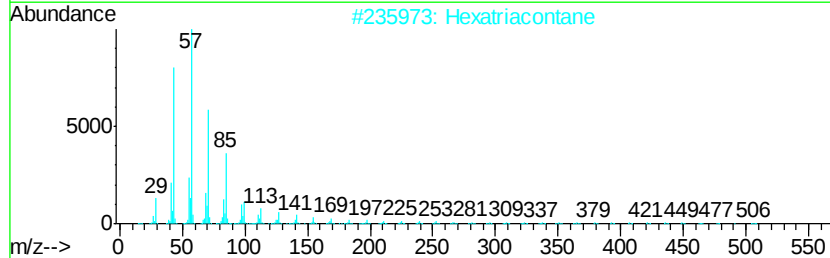
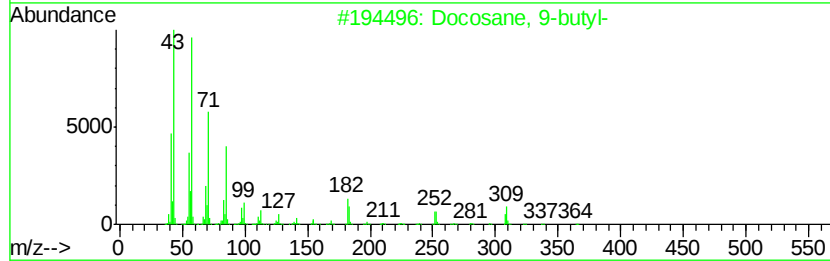
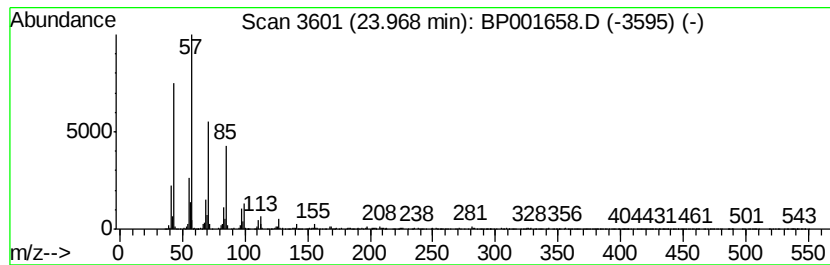
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Docosane, 9-butyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.97	12.16 ng	244916	Perylene-d12	23.37

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Docosane, 9-butyl-	366	C26H54	055282-14-9	94
2		Hexatriacontane	507	C36H74	000630-06-8	91
3		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
4		Hentriacontane	437	C31H64	000630-04-6	89
5		Docosane, 11-butyl-	366	C26H54	013475-76-8	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP011720\
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Misc :
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Instrument :
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ClientSampleId :
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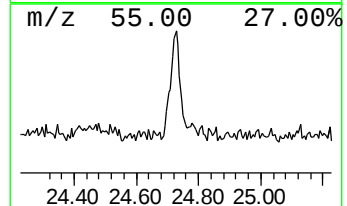
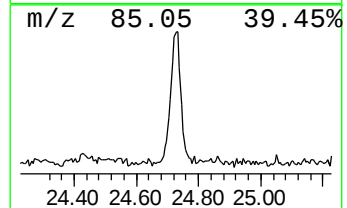
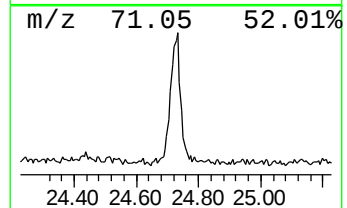
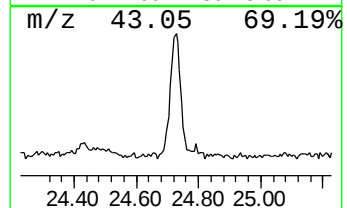
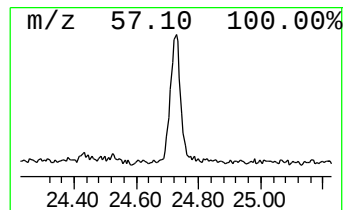
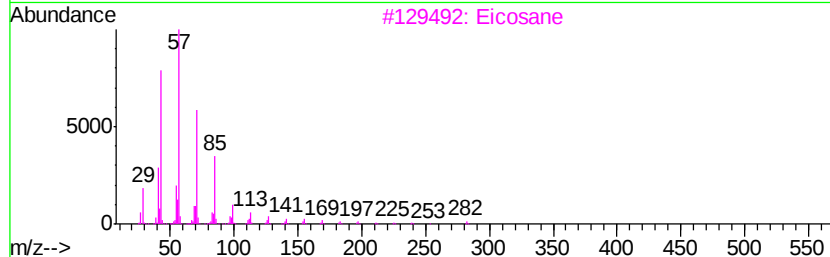
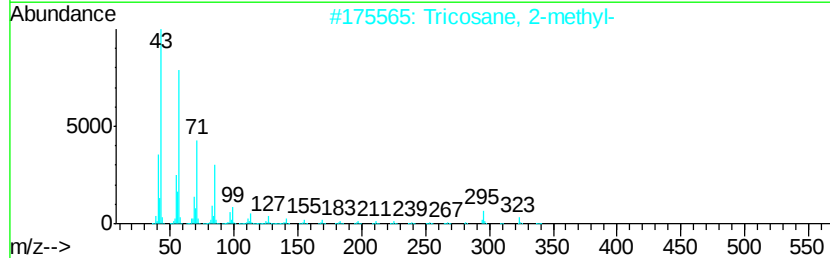
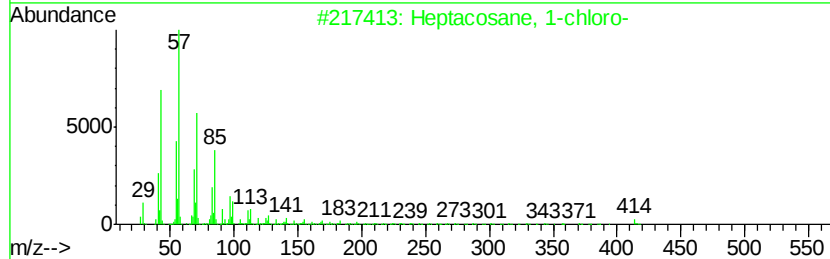
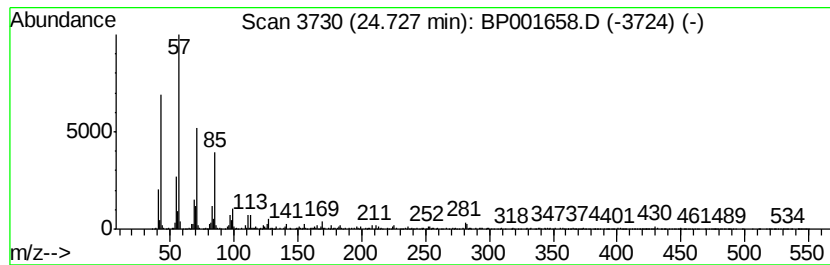
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Heptacosane, 1-chloro- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.73	6.63 ng	133536	Perylene-d12	23.37

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptacosane, 1-chloro-	414	C27H55Cl	062016-79-9	91
2		Tricosane, 2-methyl-	338	C24H50	001928-30-9	91
3		Eicosane	282	C20H42	000112-95-8	91
4		Tridecane, 7-hexyl-	268	C19H40	007225-66-3	87
5		Nonadecane	268	C19H40	000629-92-5	83



Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP011720\
 Data File : BP001658.D
 Acq On : 17 Jan 2020 15:39
 Operator : JU
 Sample : L1171-01
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BW-1A-20.0

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP010820.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
Butane, 2-methoxy...	2.95	17.8	ng	283005	1	7.63	317541	20.0
2-Pentanone, 4-hy...	4.79	15.9	ng	252254	1	7.63	317541	20.0
unknown7.18	7.18	81.5	ng	1293910	1	7.63	317541	20.0
Cobalt, [(1,2,5,6...	19.82	5.0	ng	142071	5	21.15	567903	20.0
1-Docosene	20.89	10.2	ng	288082	5	21.15	567903	20.0
Hexacosane	21.32	5.4	ng	153054	5	21.15	567903	20.0
Tridecane, 7-hexyl-	21.76	7.7	ng	219897	5	21.15	567903	20.0
Tetracosane	22.22	14.7	ng	415932	5	21.15	567903	20.0
Docosane, 11-butyl-	22.74	19.0	ng	381946	6	23.37	402712	20.0
Hentriacontane	23.31	15.9	ng	320833	6	23.37	402712	20.0
Docosane, 9-butyl-	23.97	12.2	ng	244916	6	23.37	402712	20.0
Heptacosane, 1-ch...	24.73	6.6	ng	133536	6	23.37	402712	20.0