

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP011620\
 Data File : BP001632.D
 Acq On : 16 Jan 2020 17:35
 Operator : JU
 Sample : L1145-02
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 MH-109

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	2.870	9	14	22	rBV	53646	95513	3.76%	0.555%
2	2.946	22	27	42	rVB	568731	765423	30.14%	4.451%
3	4.793	336	341	350	rVB	64801	91495	3.60%	0.532%
4	5.258	414	420	428	rBV	440926	633050	24.93%	3.681%
5	6.822	679	686	697	rBV	286845	470738	18.54%	2.737%
6	7.169	738	745	762	rBV	777461	1261020	49.66%	7.332%
7	7.628	817	823	833	rVB	165232	254018	10.00%	1.477%
8	7.946	869	877	888	rBV	650700	1056530	41.60%	6.143%
9	8.775	1010	1018	1034	rBV	543369	926722	36.49%	5.389%
10	10.405	1286	1295	1308	rBV	173098	304778	12.00%	1.772%
11	12.899	1711	1719	1733	rBV	1232790	1850613	72.87%	10.761%
12	13.740	1856	1862	1873	rBV	81453	116707	4.60%	0.679%
13	14.275	1947	1953	1962	rBV2	268036	405112	15.95%	2.356%
14	15.781	2202	2209	2216	rBV	1067748	1550069	61.04%	9.013%
15	15.840	2216	2219	2231	rVB8	14980	30415	1.20%	0.177%
16	17.040	2417	2423	2433	rBV	339171	477494	18.80%	2.776%
17	17.975	2577	2582	2594	rBV2	108357	190455	7.50%	1.107%
18	19.628	2857	2863	2870	rBV	2154536	2539499	100.00%	14.766%
19	20.439	2998	3001	3005	rVB	42312	45393	1.79%	0.264%
20	20.892	3073	3078	3084	rBV	285843	407549	16.05%	2.370%
21	21.145	3116	3121	3128	rVV	328860	443734	17.47%	2.580%
22	21.322	3148	3151	3155	rVB	164057	175842	6.92%	1.022%
23	21.763	3221	3226	3238	rVB	243562	331903	13.07%	1.930%
24	22.227	3301	3305	3311	rVB	314767	400451	15.77%	2.328%
25	22.357	3322	3327	3334	rVB	313012	431496	16.99%	2.509%
26	22.739	3387	3392	3401	rVB	311030	442461	17.42%	2.573%
27	23.316	3485	3490	3495	rBV	266435	437095	17.21%	2.542%
28	23.375	3495	3500	3511	rVB	215583	403743	15.90%	2.348%
29	23.974	3596	3602	3611	rBV	200677	391102	15.40%	2.274%
30	24.733	3727	3731	3738	rVB	98407	197831	7.79%	1.150%
31	25.621	3878	3882	3890	rVB3	31148	69796	2.75%	0.406%

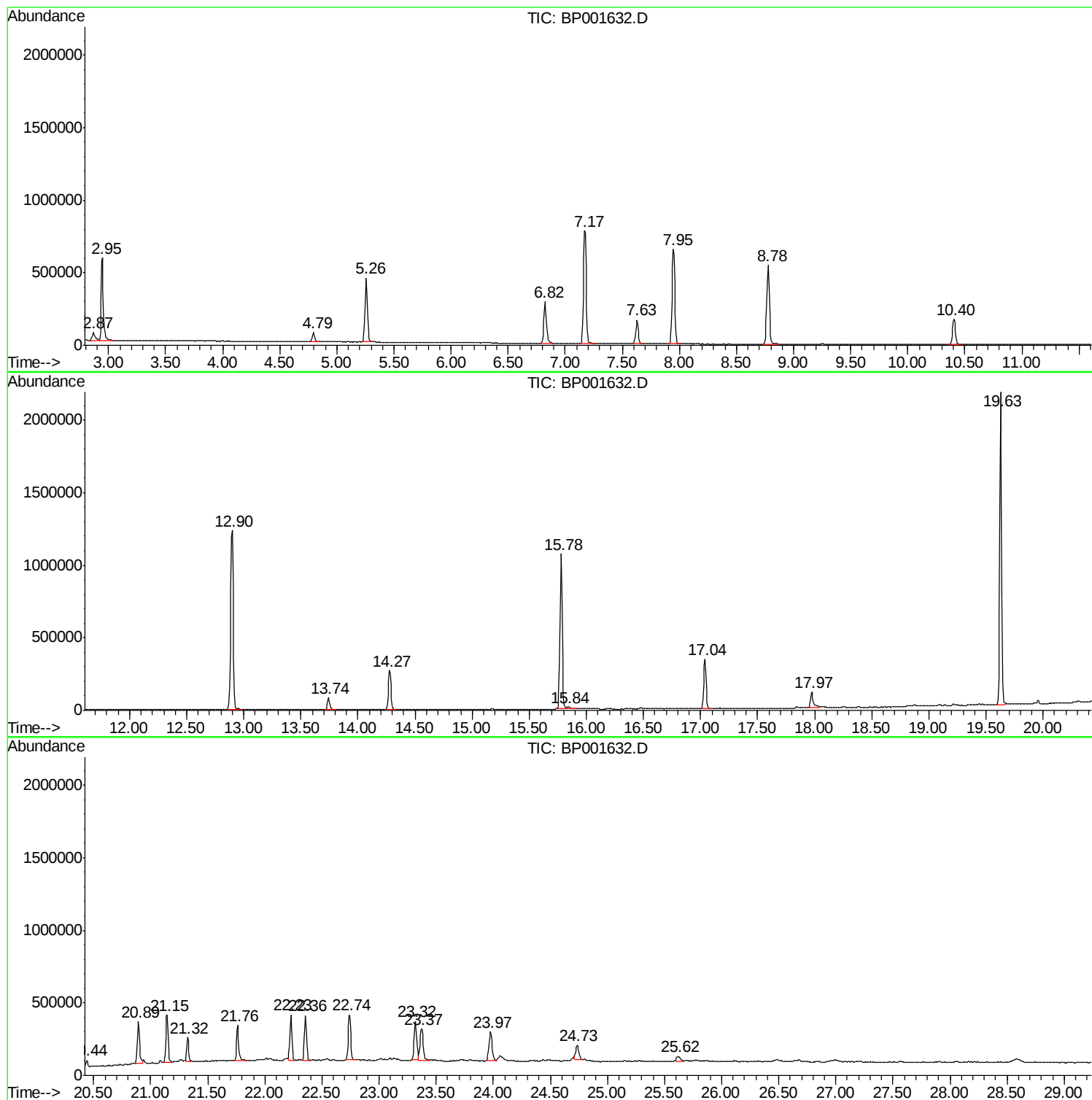
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP010820.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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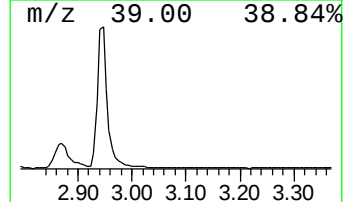
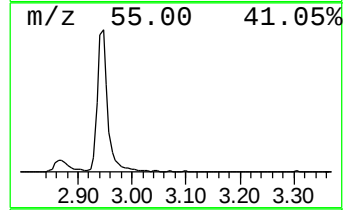
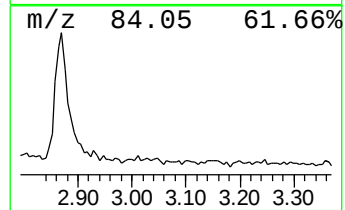
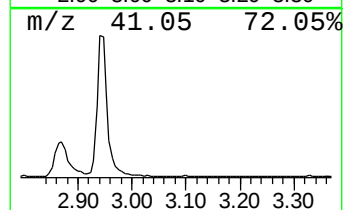
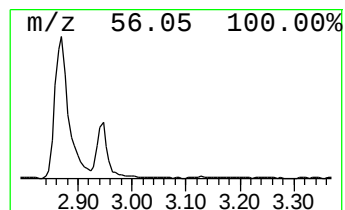
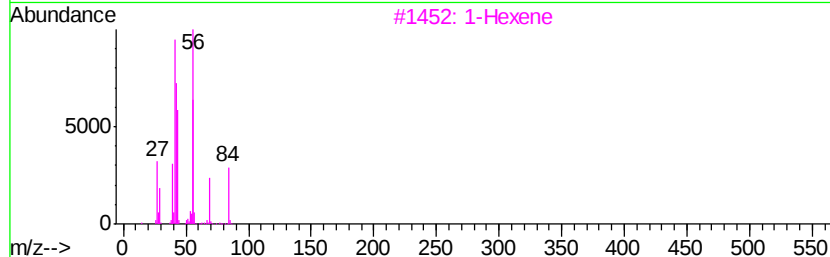
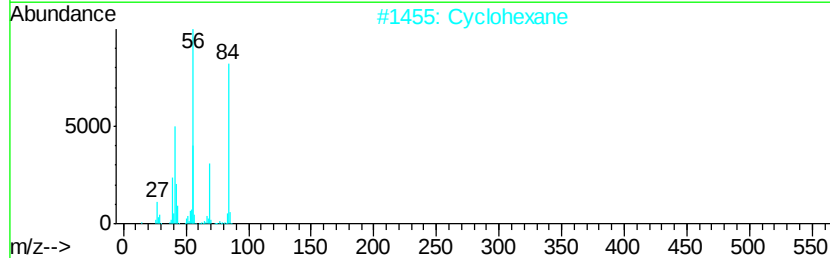
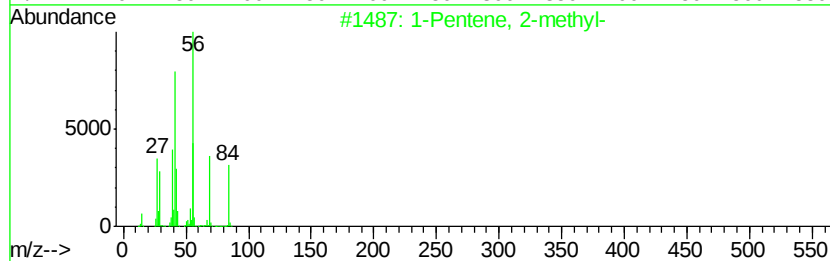
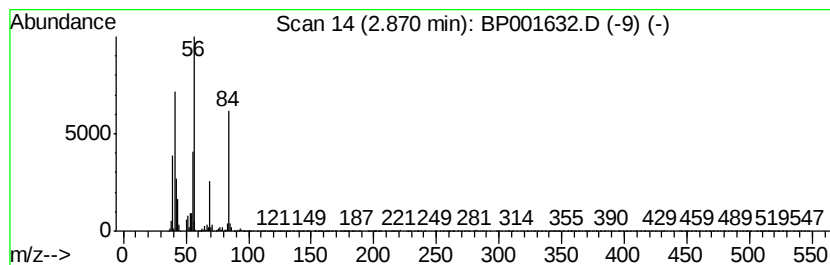
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 1-Pentene, 2-methyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.87	7.52 ng	95513	1,4-Dichlorobenzene-d4	7.63

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Pentene, 2-methyl-	84	C6H12	000763-29-1	86
2			Cyclohexane	84	C6H12	000110-82-7	81
3			1-Hexene	84	C6H12	000592-41-6	47
4			Cyanoic acid, 2-methylpropyl ester	99	C5H9NO	001768-25-8	45
5			Cyclobutane, ethyl-	84	C6H12	004806-61-5	43



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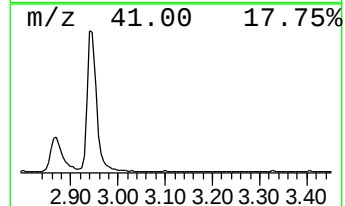
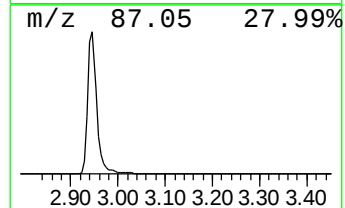
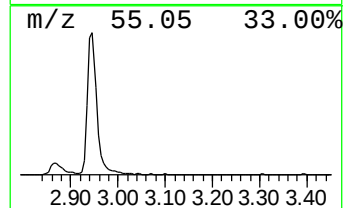
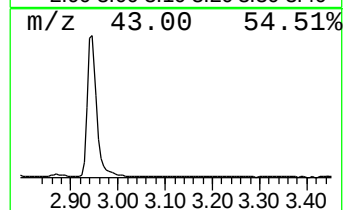
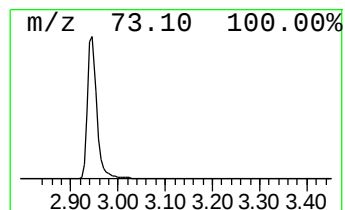
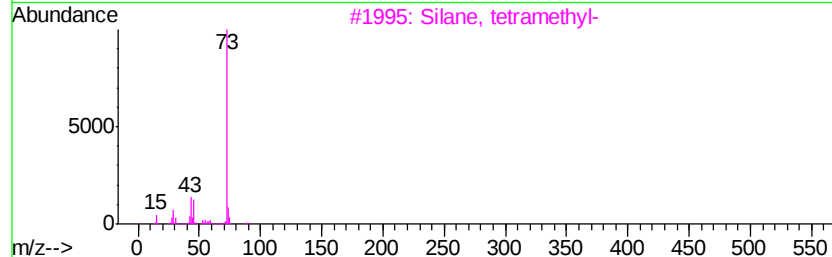
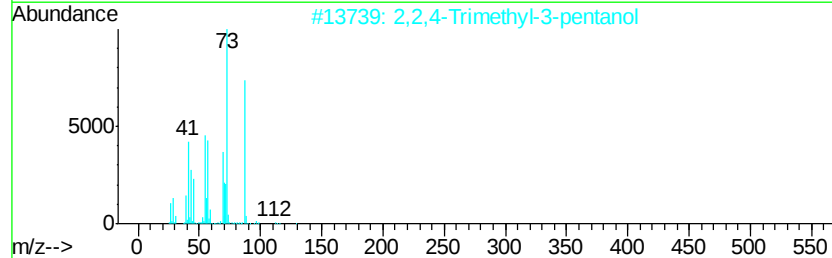
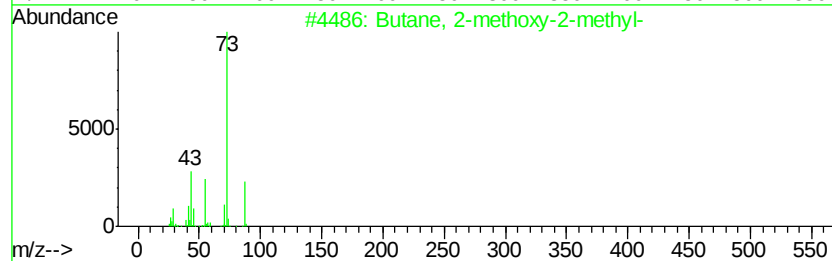
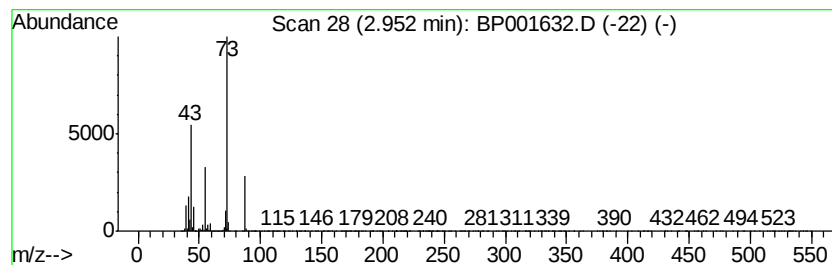
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.95	60.27 ng	765423	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	90
2		2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	40
3		Silane, tetramethyl-	88	C4H12Si	000075-76-3	32
4		2-Methyl-5-hexen-3-ol	114	C7H14O	032815-70-6	23
5		1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	9



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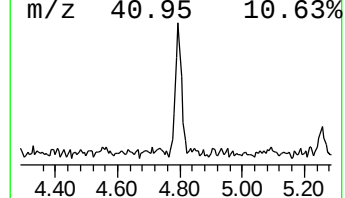
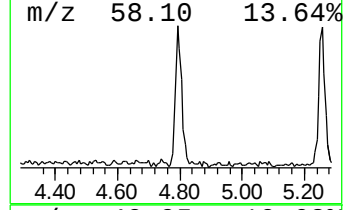
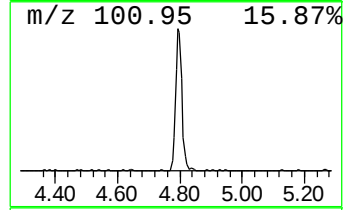
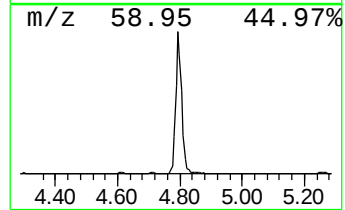
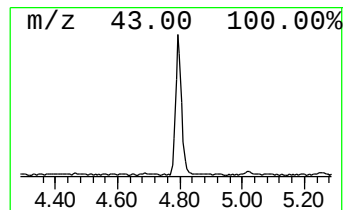
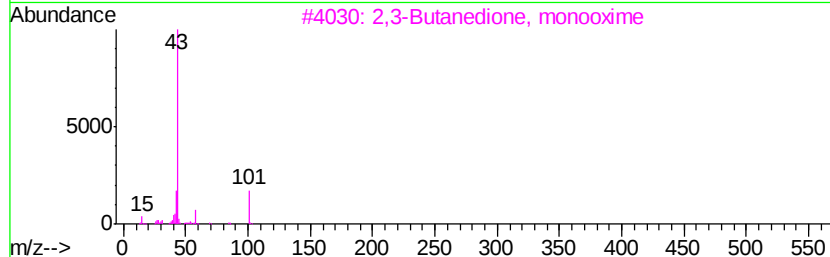
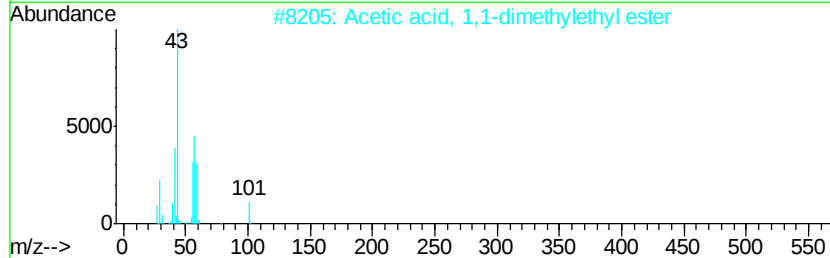
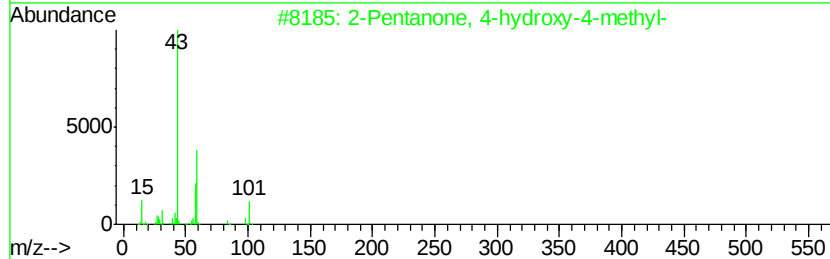
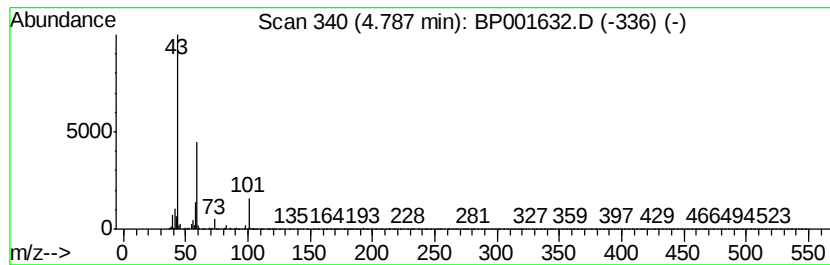
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TIC Integration Parameters: LSCINT.P

Peak Number 3 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.79	7.20 ng	91495	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	50
2		Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	28
3		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	16
4		1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10
5		(+)-4-Amino-4,5-dihydro-2(3H)-f...	101	C4H7NO2	016504-58-8	9



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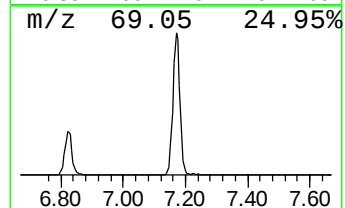
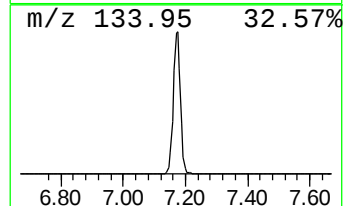
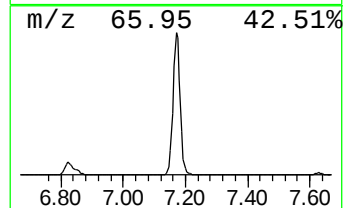
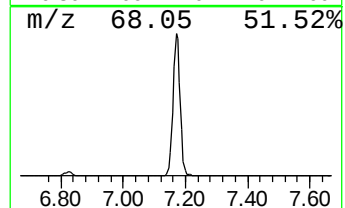
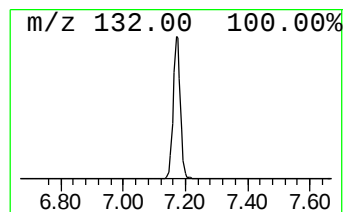
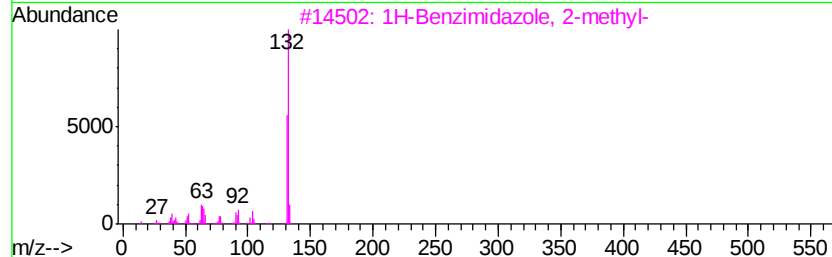
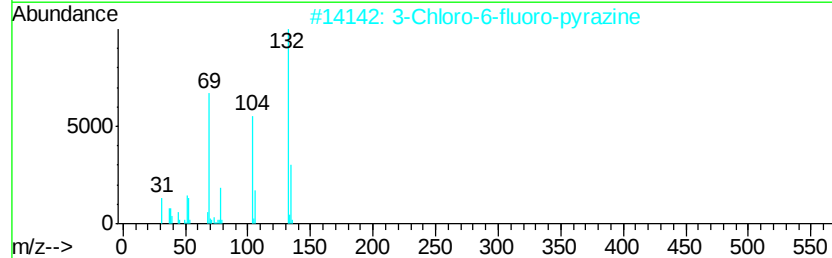
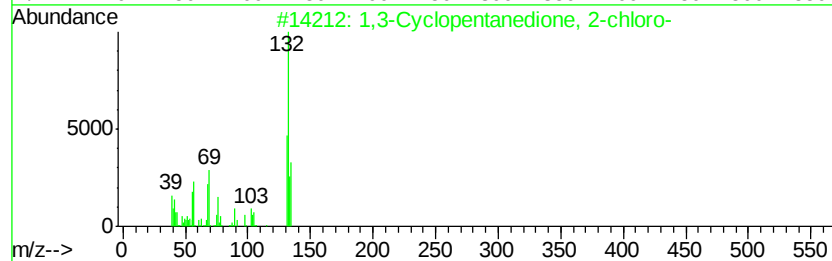
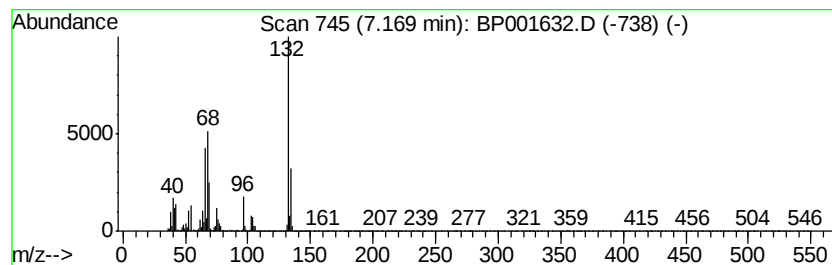
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 4 unknown7.17 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.17	99.29 ng	1261020	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	27
2		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	14
3		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	14
4		3H-1,2-Dithiol-3-one, 5-methyl-	132	C4H4OS2	003620-08-4	14
5		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10



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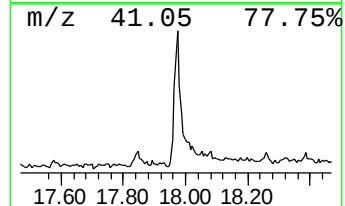
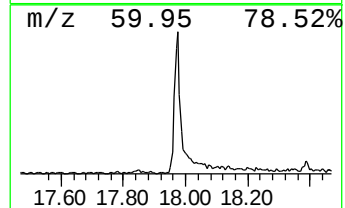
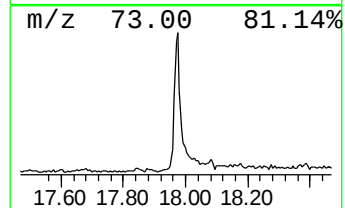
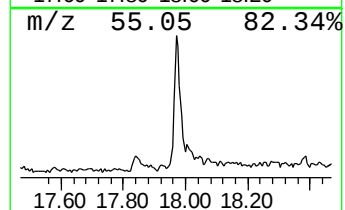
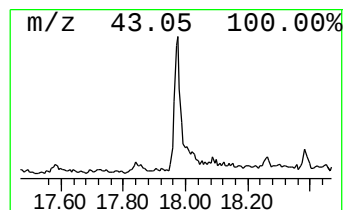
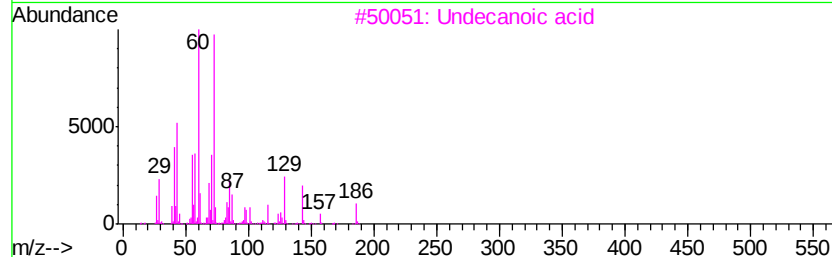
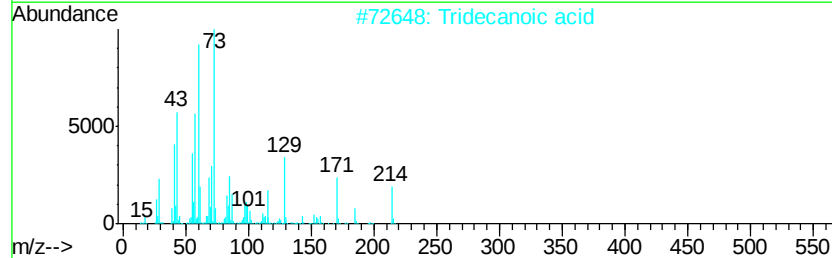
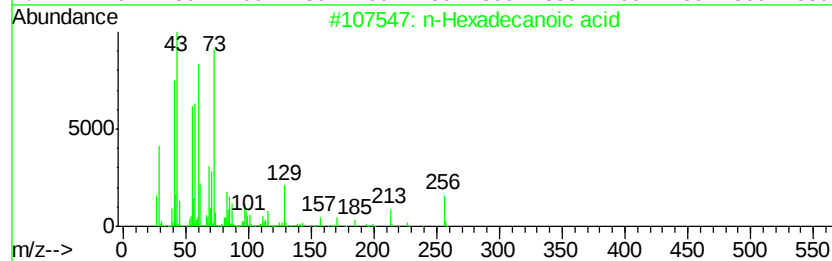
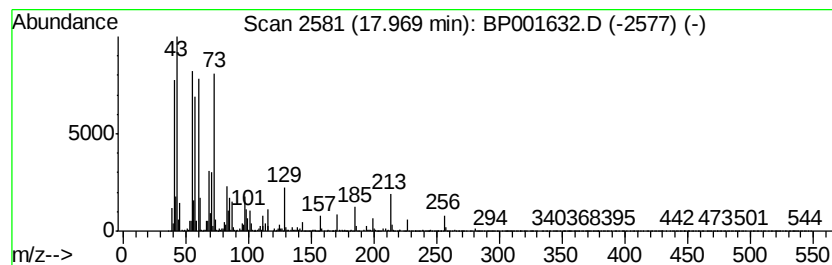
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TIC Library : C:\DATABASE\NIST11.L
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 Peak Number 6 n-Hexadecanoic acid Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.97	7.98 ng	190455	Phenanthrene-d10	17.04

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	96
2		Tridecanoic acid	214	C13H26O2	000638-53-9	95
3		Undecanoic acid	186	C11H22O2	000112-37-8	70
4		Octadecanoic acid	284	C18H36O2	000057-11-4	70
5		Isopropyl palmitate	298	C19H38O2	000142-91-6	64



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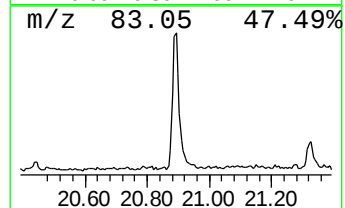
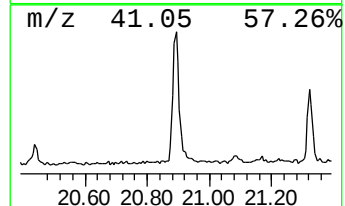
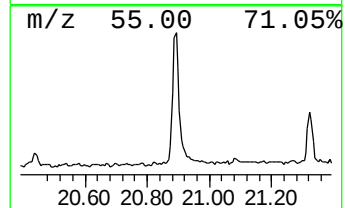
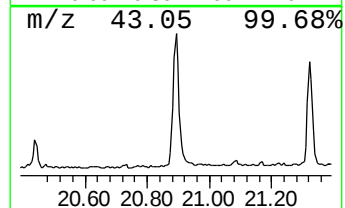
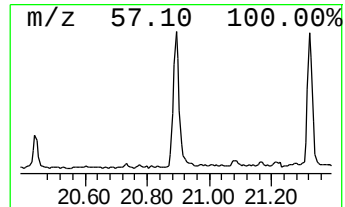
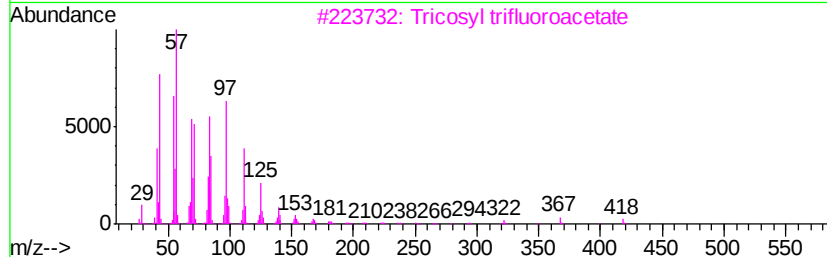
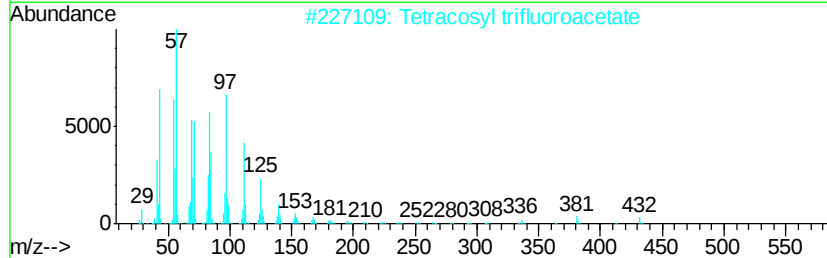
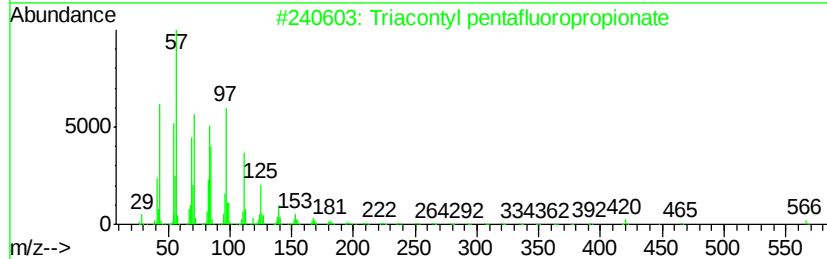
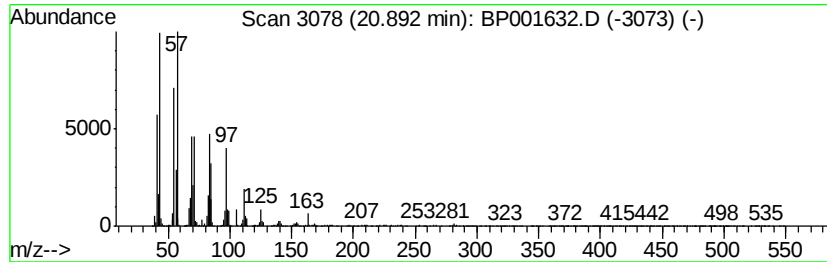
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Triacontyl pentafluoropropi... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.89	18.37 ng	407549	Chrysene-d12	21.15

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Triacontyl pentafluoropropionate	584	C33H61F5O2	1000351-80-0	91
2		Tetracosyl trifluoroacetate	450	C26H49F3O2	1000351-76-4	91
3		Tricosyl trifluoroacetate	436	C25H47F3O2	1000351-75-1	91
4		Hexacosyl pentafluoropropionate	528	C29H53F5O2	1000351-81-2	91
5		Hexacosyl heptafluorobutyrate	578	C30H53F7O2	1000351-83-3	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP011620\
Data File : BP001632.D
Acq On : 16 Jan 2020 17:35
Operator : JU
Sample : L1145-02
Misc :
ALS Vial : 4 Sample Multiplier: 1

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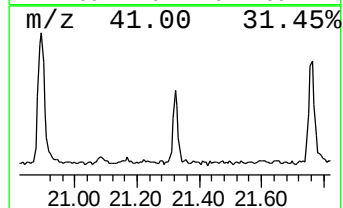
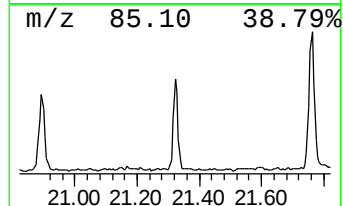
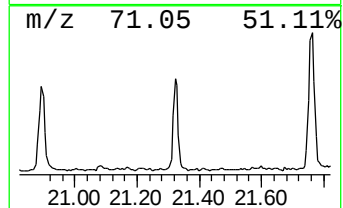
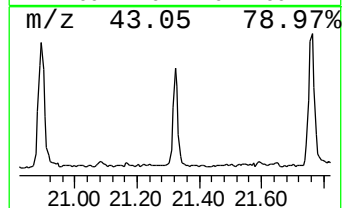
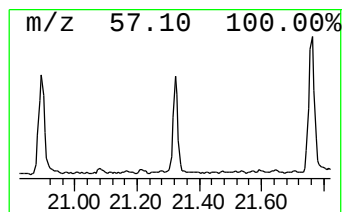
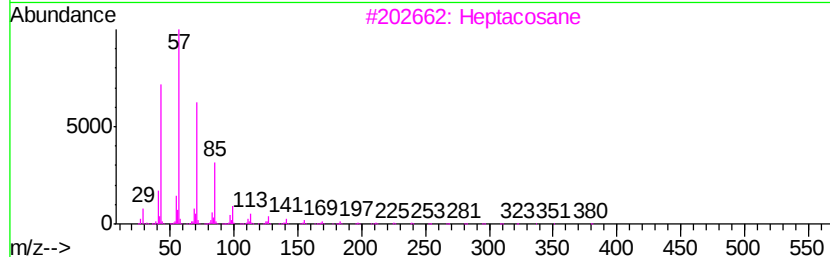
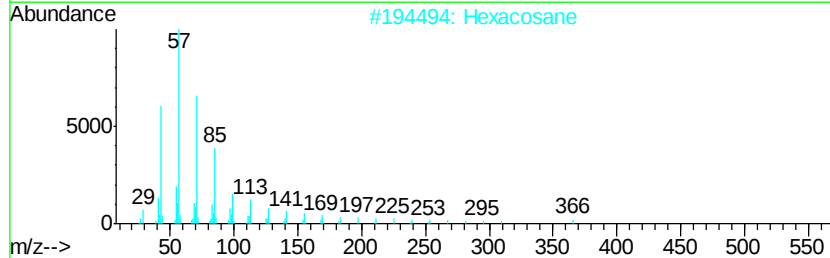
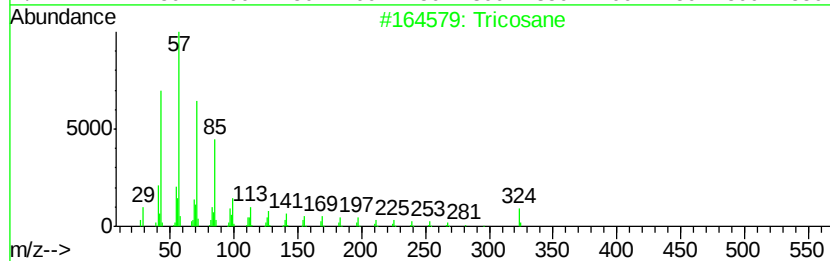
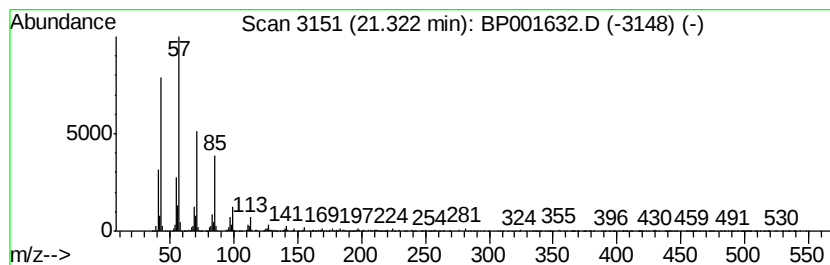
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Tricosane Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.32	7.93 ng	175842	Chrysene-d12	21.15

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tricosane		324	C23H48	000638-67-5	93
2	Hexacosane		366	C26H54	000630-01-3	92
3	Heptacosane		380	C27H56	000593-49-7	90
4	Hexadecane, 3-methyl-		240	C17H36	006418-43-5	87
5	Eicosane, 10-methyl-		296	C21H44	054833-23-7	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP011620\
Data File : BP001632.D
Acq On : 16 Jan 2020 17:35
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Misc :
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ClientSampleId :
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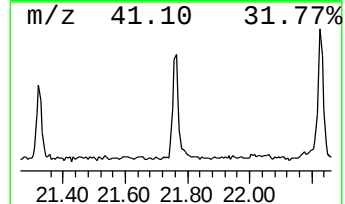
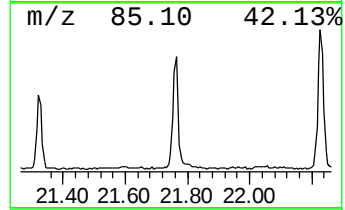
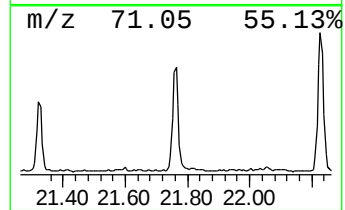
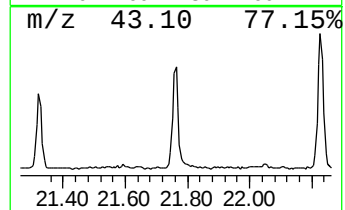
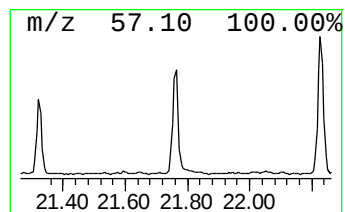
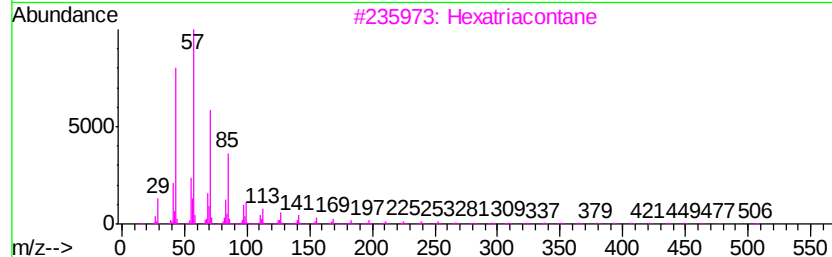
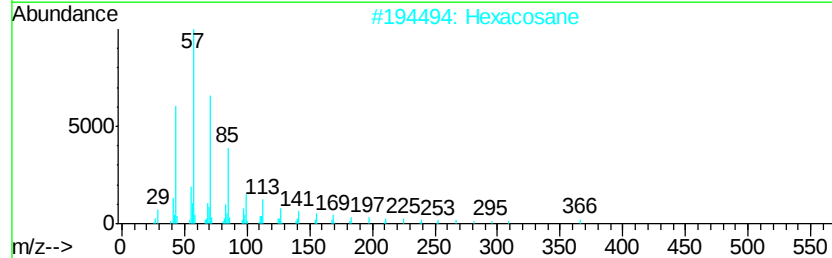
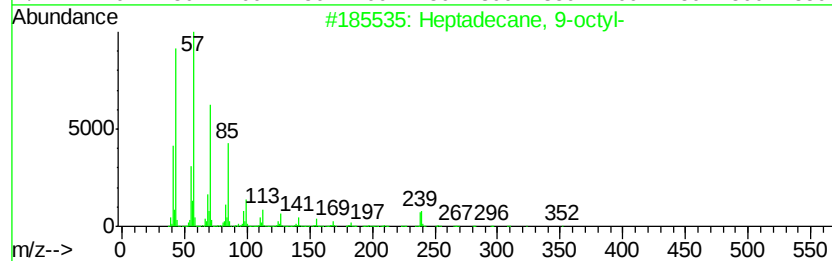
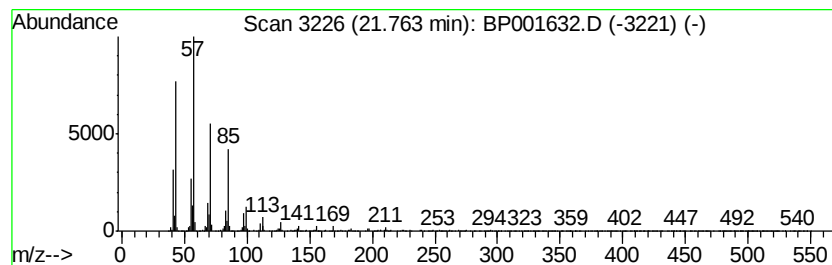
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Heptadecane, 9-octyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.76	14.96 ng	331903	Chrysene-d12	21.15

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	95
2		Hexacosane	366	C26H54	000630-01-3	94
3		Hexatriacontane	507	C36H74	000630-06-8	93
4		Pentacosane	352	C25H52	000629-99-2	93
5		Tridecane, 1-iodo-	310	C13H27I	035599-77-0	93



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP011620\
Data File : BP001632.D
Acq On : 16 Jan 2020 17:35
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Sample : L1145-02
Misc :
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Instrument :
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ClientSampleId :
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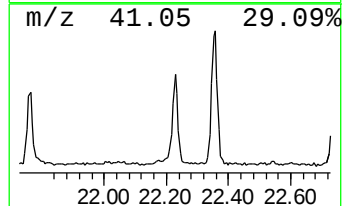
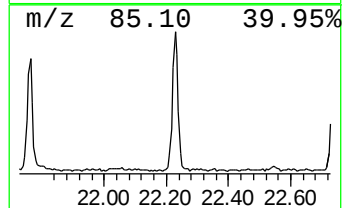
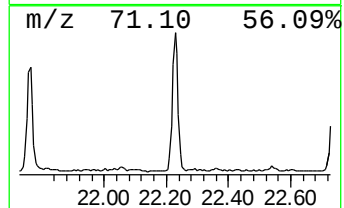
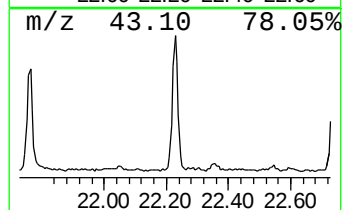
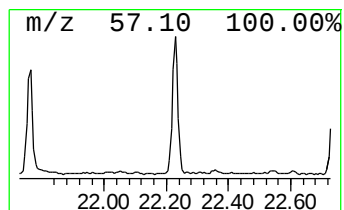
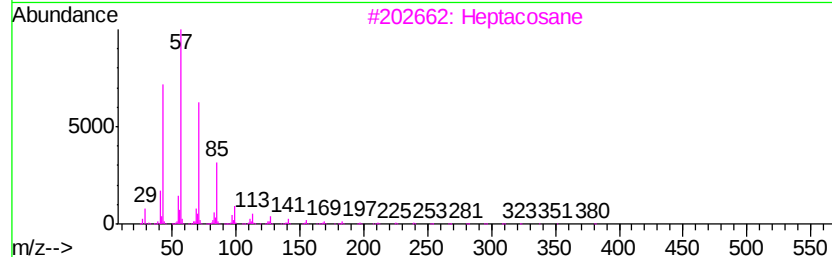
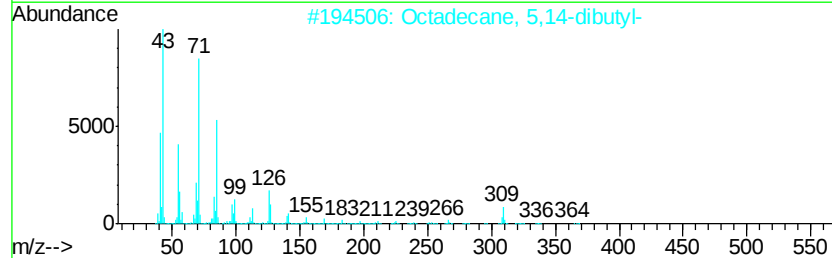
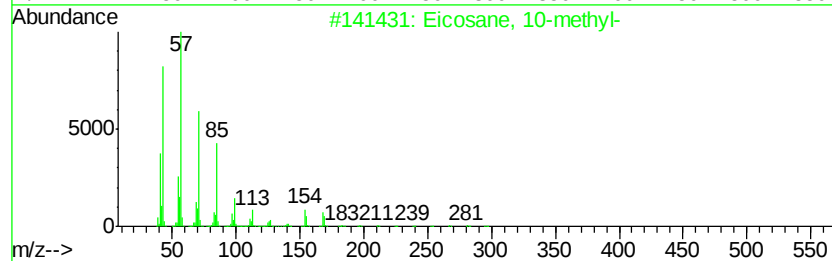
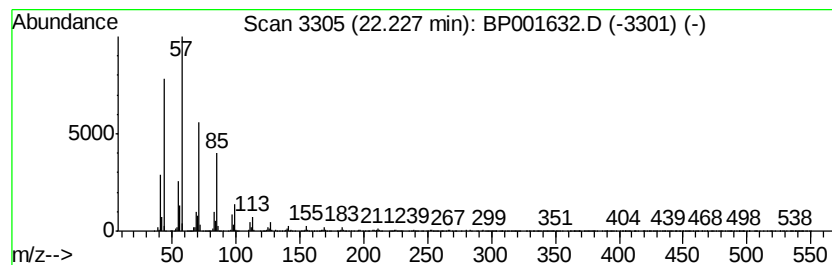
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Eicosane, 10-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.23	18.05 ng	400451	Chrysene-d12	21.15

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane, 10-methyl-	296	C21H44	054833-23-7	94
2		Octadecane, 5,14-dibutyl-	366	C26H54	055282-13-8	93
3		Heptacosane	380	C27H56	000593-49-7	91
4		Tridecane, 7-propyl-	226	C16H34	055045-09-5	91
5		1-Iodo-2-methylundecane	296	C12H25I	073105-67-6	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP011620\
Data File : BP001632.D
Acq On : 16 Jan 2020 17:35
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Misc :
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ClientSampleId :
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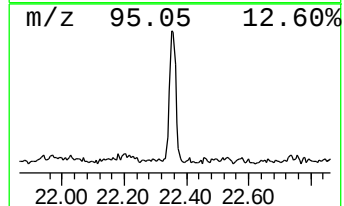
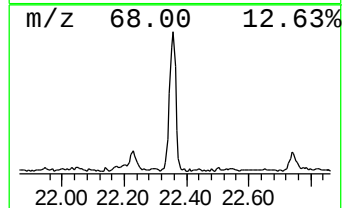
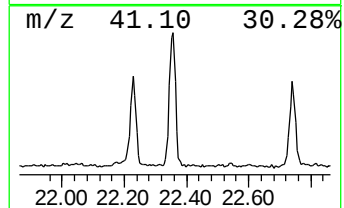
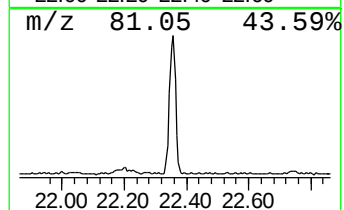
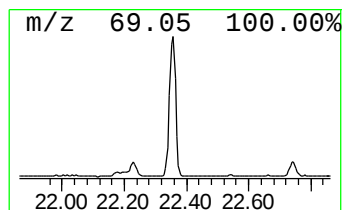
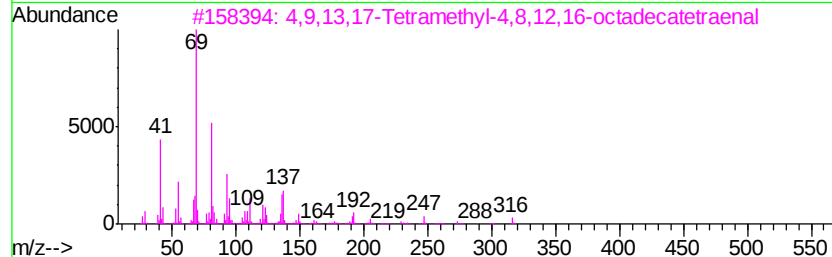
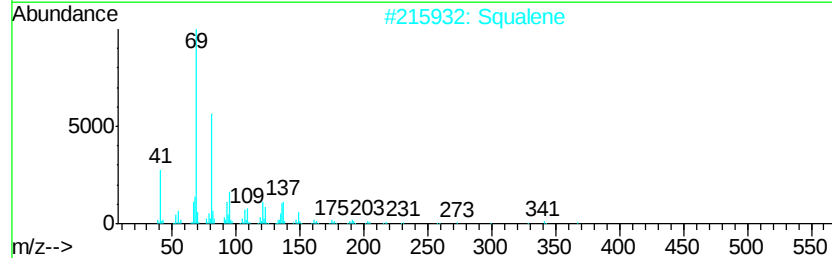
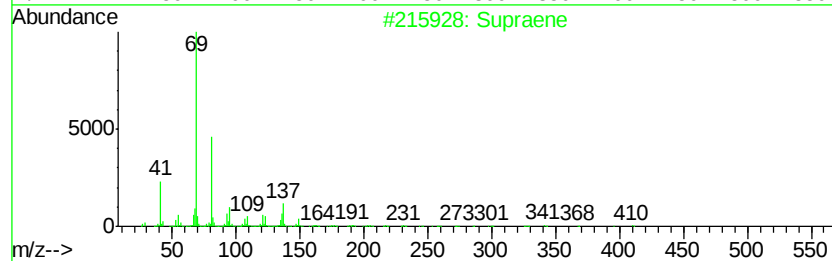
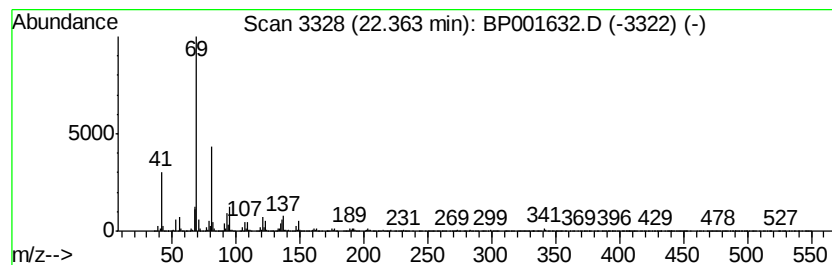
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Supraene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.36	21.37 ng	431496	Perylene-d12	23.37

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Supraene	410	C30H50	007683-64-9	91
2		Squalene	410	C30H50	000111-02-4	90
3		4,9,13,17-Tetramethyl-4,8,12,16-...	316	C22H36O	056882-09-8	83
4		4,8,12-Tetradecatrienal, 5,9,13-...	248	C17H28O	066408-55-7	72
5		Bis(3-methylbut-3-enyl) phthalate	302	C18H22O4	022711-53-1	59



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP011620\
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Sample : L1145-02
Misc :
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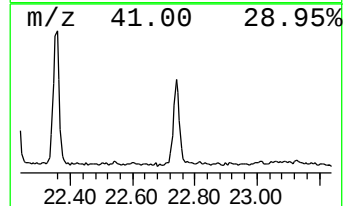
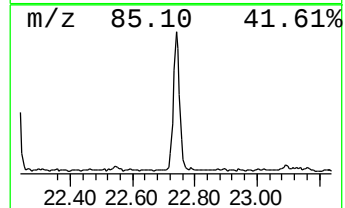
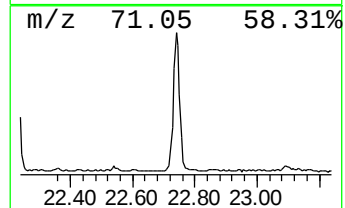
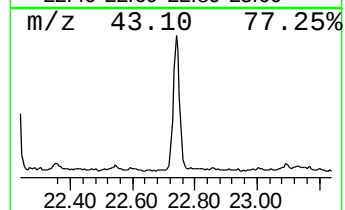
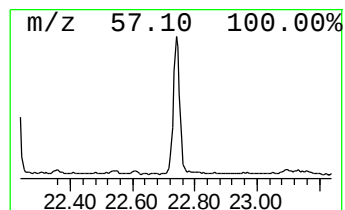
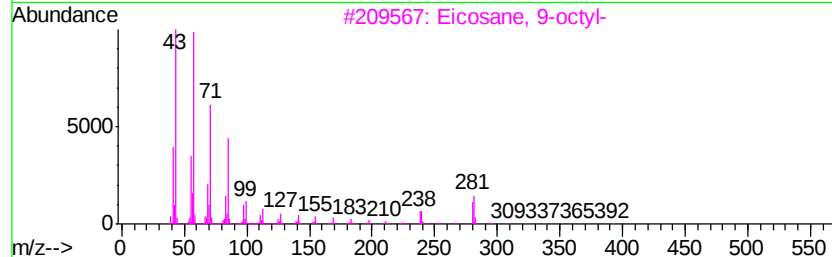
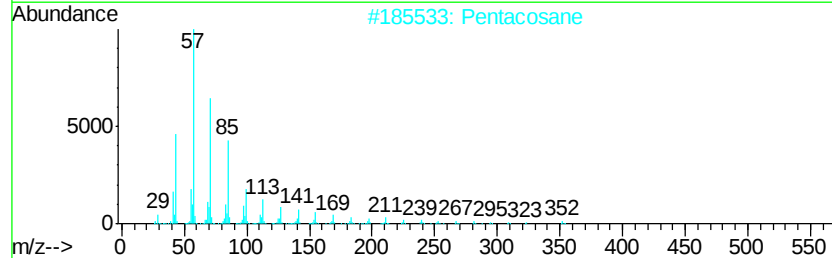
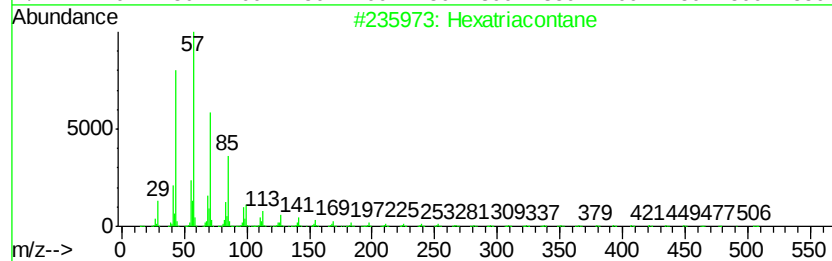
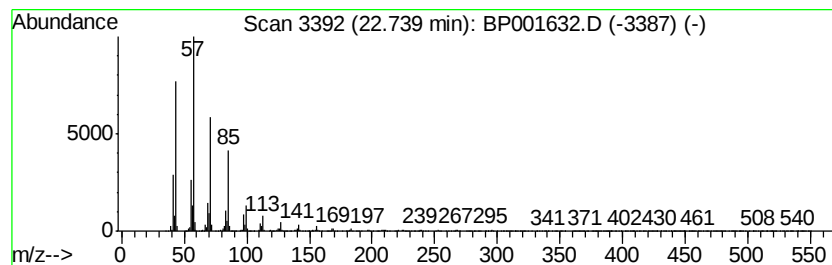
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Hexatriacontane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.74	21.92 ng	442461	Perylene-d12	23.37

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexatriacontane	507	C36H74	000630-06-8	98
2		Pentacosane	352	C25H52	000629-99-2	96
3		Eicosane, 9-octyl-	394	C28H58	013475-77-9	95
4		Docosane, 11-butyl-	366	C26H54	013475-76-8	94
5		Heneicosane, 11-decyl-	437	C31H64	055320-06-4	94



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP011620\
 Data File : BP001632.D
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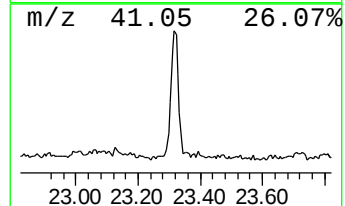
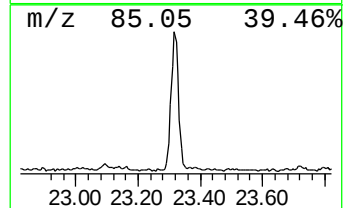
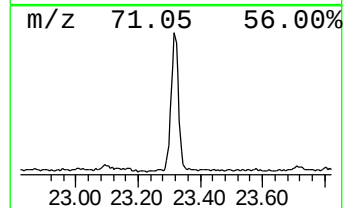
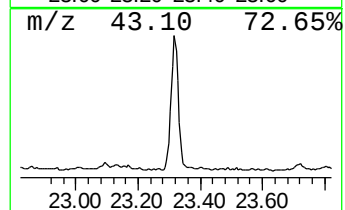
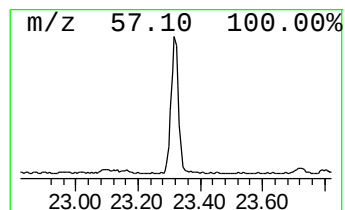
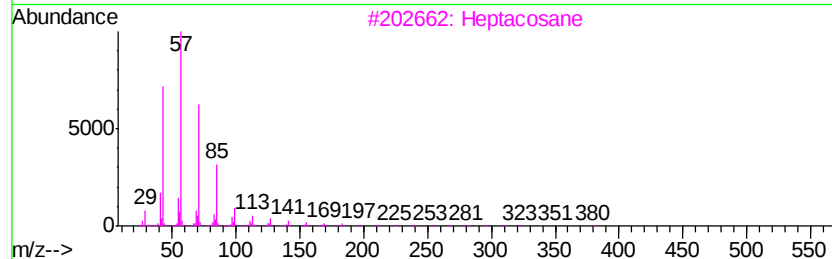
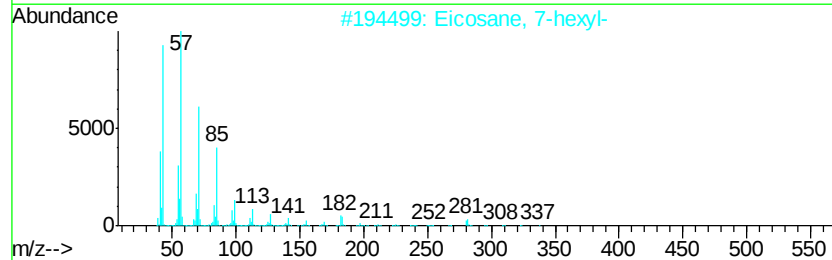
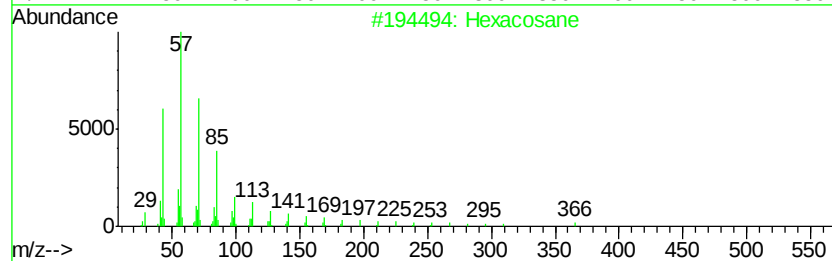
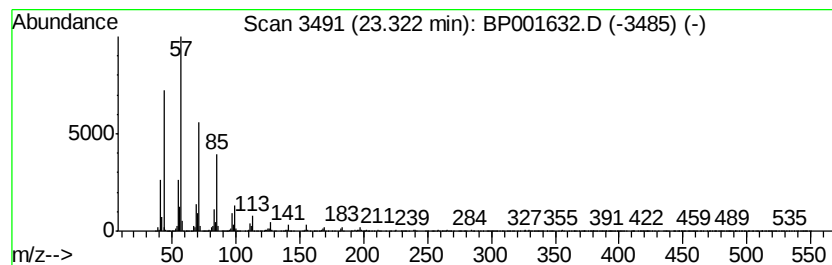
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TIC Library : C:\DATABASE\NIST11.L
 TIC Integration Parameters: LSCINT.P

 Peak Number 14 Hexacosane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.32	21.65 ng	437095	Perylene-d12	23.37

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexacosane		366	C26H54	000630-01-3	94
2	Eicosane, 7-hexyl-		366	C26H54	055333-99-8	91
3	Heptacosane		380	C27H56	000593-49-7	91
4	Tetratetracontane		619	C44H90	007098-22-8	91
5	Heptadecane, 9-octyl-		352	C25H52	007225-64-1	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP011620\
Data File : BP001632.D
Acq On : 16 Jan 2020 17:35
Operator : JU
Sample : L1145-02
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
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ClientSampleId :
MH-109

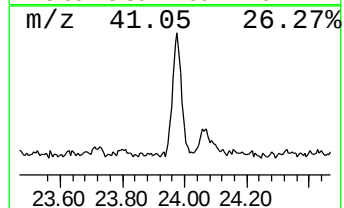
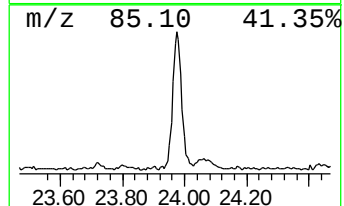
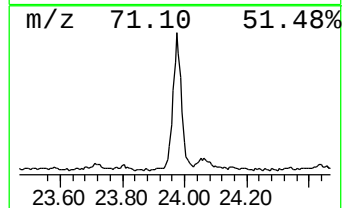
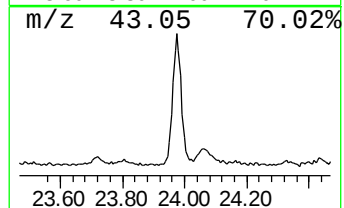
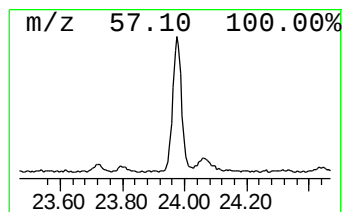
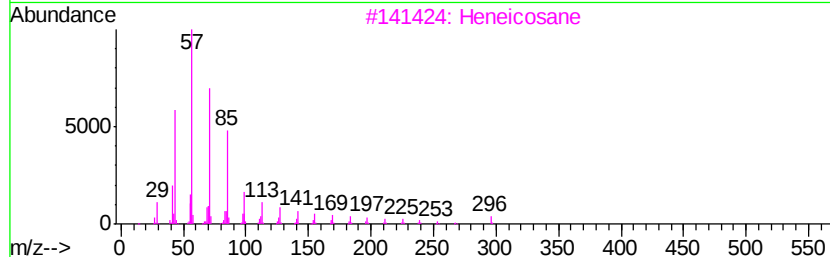
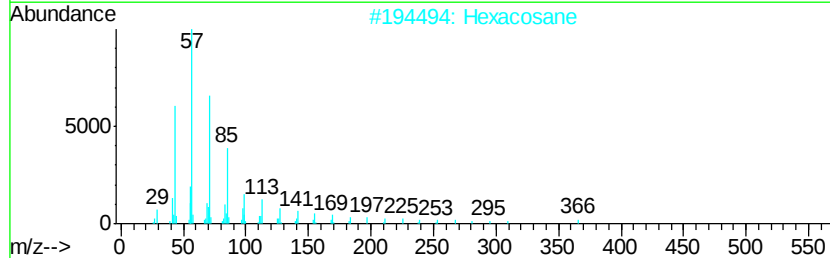
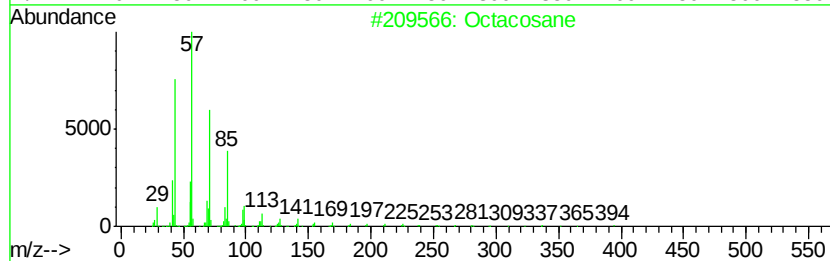
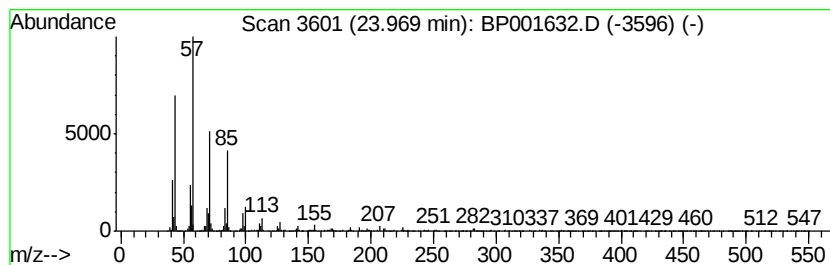
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Octacosane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.97	19.37 ng	391102	Perylene-d12	23.37

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octacosane	394	C28H58	000630-02-4	98
2		Hexacosane	366	C26H54	000630-01-3	96
3		Heneicosane	296	C21H44	000629-94-7	95
4		Eicosane	282	C20H42	000112-95-8	95
5		Eicosane, 9-octyl-	394	C28H58	013475-77-9	95



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP011620\
 Data File : BP001632.D
 Acq On : 16 Jan 2020 17:35
 Operator : JU
 Sample : L1145-02
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 MH-109

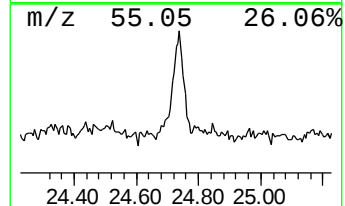
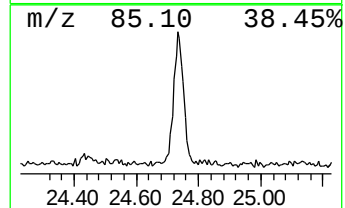
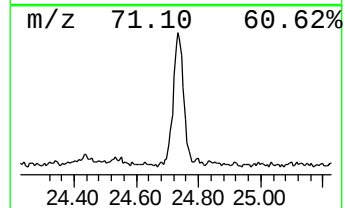
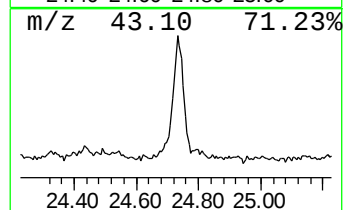
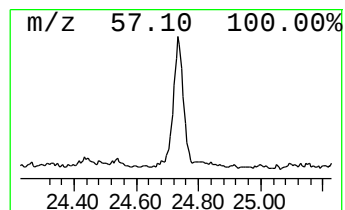
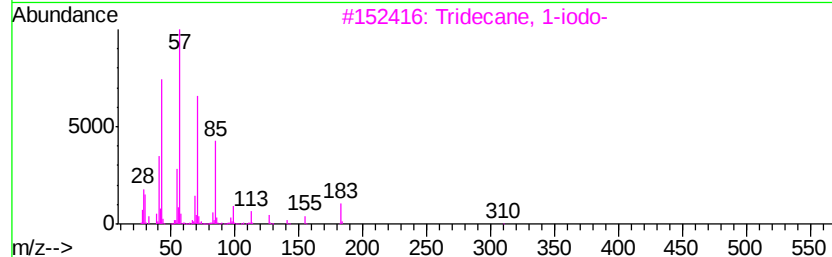
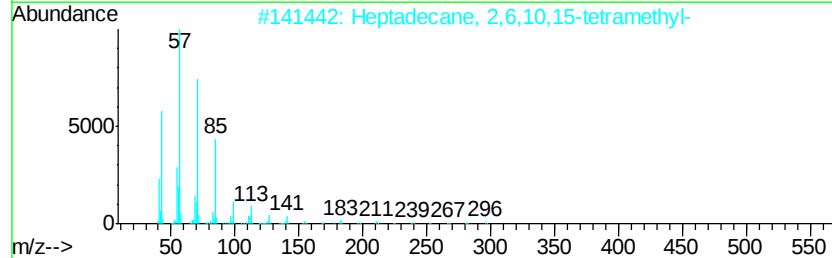
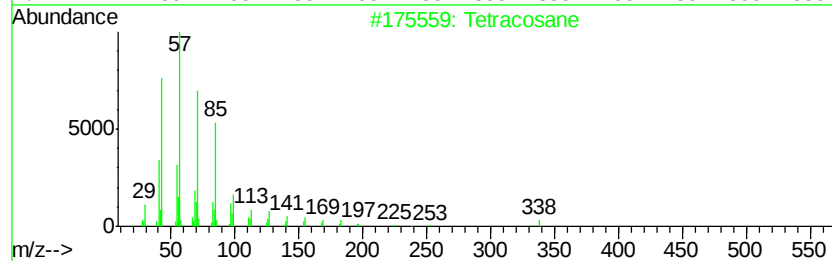
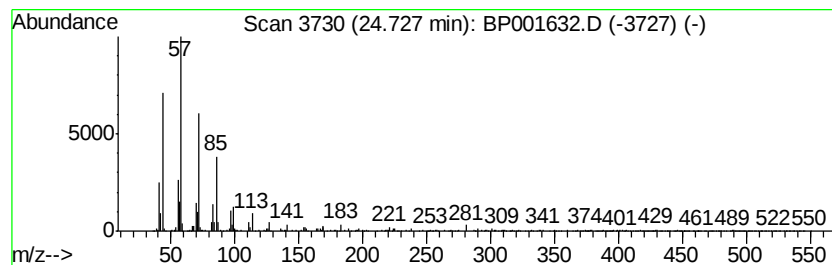
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

 Peak Number 16 Tetracosane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.73	9.80 ng	197831	Perylene-d12	23.37

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetracosane	338	C24H50	000646-31-1	93
2		Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	93
3		Tridecane, 1-iodo-	310	C13H27I	035599-77-0	93
4		Hexadecane, 6,11-dipentyl-	366	C26H54	015874-03-0	90
5		Octacosane	394	C28H58	000630-02-4	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP011620\
Data File : BP001632.D
Acq On : 16 Jan 2020 17:35
Operator : JU
Sample : L1145-02
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
MH-109

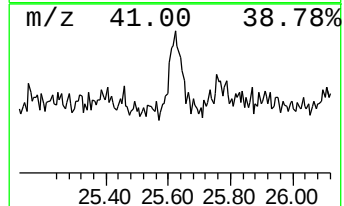
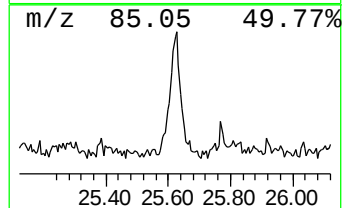
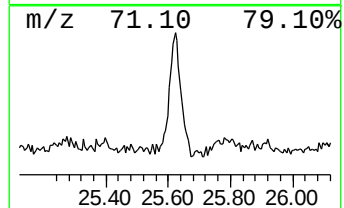
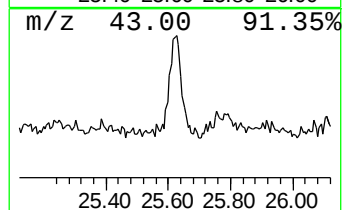
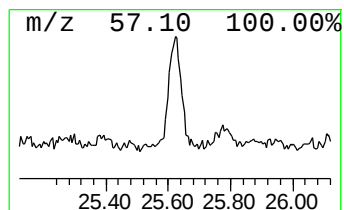
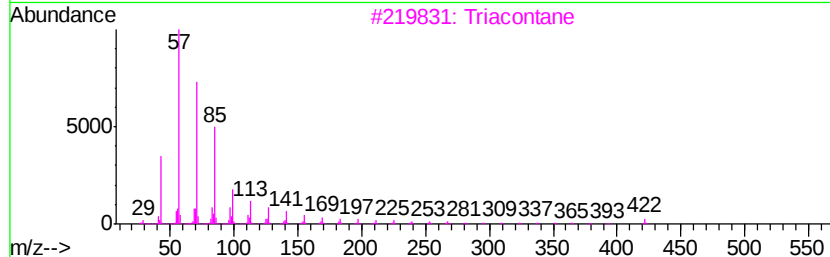
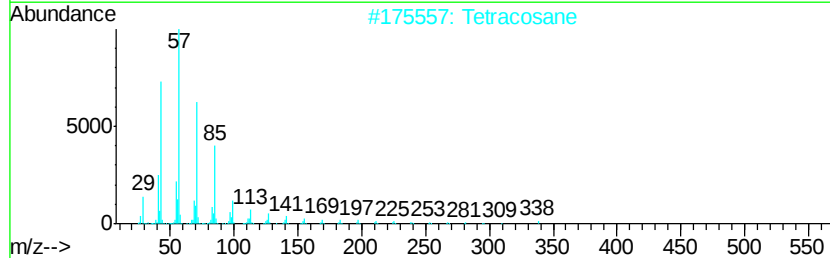
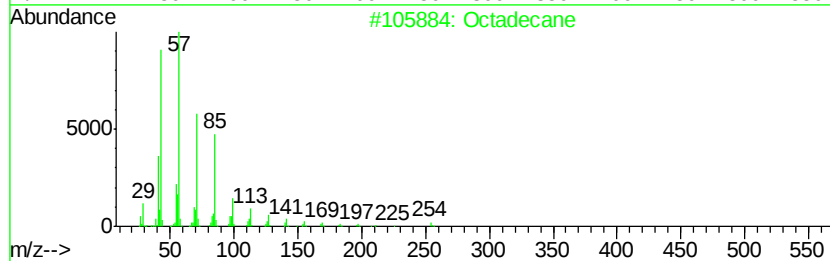
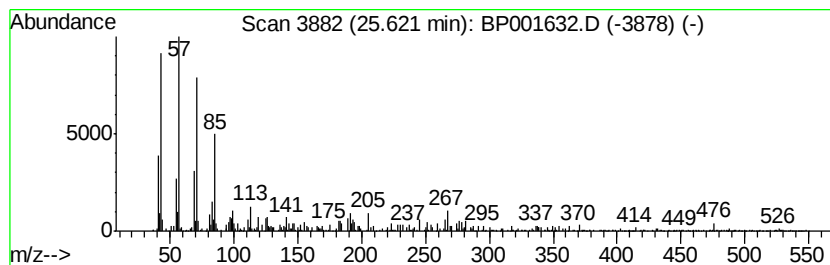
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 Octadecane Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.62	3.46 ng	69796	Perylene-d12	23.37

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecane	254	C18H38	000593-45-3	91
2			Tetracosane	338	C24H50	000646-31-1	91
3			Triacotane	422	C30H62	000638-68-6	91
4			Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
5			Octacosane	394	C28H58	000630-02-4	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP011620\
 Data File : BP001632.D
 Acq On : 16 Jan 2020 17:35
 Operator : JU
 Sample : L1145-02
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 MH-109

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
1-Pentene, 2-methyl-	2.87	7.5	ng	95513	1	7.63	254018	20.0
Butane, 2-methoxy...	2.95	60.3	ng	765423	1	7.63	254018	20.0
2-Pentanone, 4-hy...	4.79	7.2	ng	91495	1	7.63	254018	20.0
unknown7.17	7.17	99.3	ng	1261020	1	7.63	254018	20.0
n-Hexadecanoic acid	17.97	8.0	ng	190455	4	17.04	477494	20.0
triacontyl penta...	20.89	18.4	ng	407549	5	21.15	443734	20.0
Tricosane	21.32	7.9	ng	175842	5	21.15	443734	20.0
Heptadecane, 9-oc...	21.76	15.0	ng	331903	5	21.15	443734	20.0
Eicosane, 10-methyl-	22.23	18.1	ng	400451	5	21.15	443734	20.0
Supraene	22.36	21.4	ng	431496	6	23.37	403743	20.0
Hexatriacontane	22.74	21.9	ng	442461	6	23.37	403743	20.0
Hexacosane	23.32	21.6	ng	437095	6	23.37	403743	20.0
Octacosane	23.97	19.4	ng	391102	6	23.37	403743	20.0
Tetracosane	24.73	9.8	ng	197831	6	23.37	403743	20.0
Octadecane	25.62	3.5	ng	69796	6	23.37	403743	20.0