

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM121918\
 Data File : BM018318.D
 Acq On : 20 Dec 2018 04:14
 Operator : JU/SJ
 Sample : J6378-24
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 P001-SS002-1218-01

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_M\METHODS\8270-BM121518.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	4.681	300	305	311	rBV	109799	146473	2.77%	0.329%
2	5.122	373	380	397	rBV	2858303	3985381	75.24%	8.951%
3	6.692	639	647	662	rBV	2615362	4348435	82.09%	9.767%
4	7.028	696	704	718	rBV	2887289	4390798	82.89%	9.862%
5	7.481	775	781	789	rBV	504738	766642	14.47%	1.722%
6	7.798	827	835	844	rBV	2131575	3344584	63.14%	7.512%
7	8.645	971	979	991	rBV	1519416	2602889	49.14%	5.846%
8	10.251	1244	1252	1262	rBV	502104	897335	16.94%	2.015%
9	12.751	1669	1677	1687	rBV	2895979	4335703	81.85%	9.738%
10	12.921	1702	1706	1713	rBV3	37068	64791	1.22%	0.146%
11	13.615	1819	1824	1832	rBV	111905	184098	3.48%	0.413%
12	14.145	1905	1914	1922	rBV2	617923	999971	18.88%	2.246%
13	15.651	2164	2170	2183	rBV	2400383	3418893	64.54%	7.679%
14	16.904	2377	2383	2389	rBV2	719098	1096022	20.69%	2.462%
15	16.951	2389	2391	2396	rVB3	50477	65505	1.24%	0.147%
16	17.821	2534	2539	2548	rBV	365963	509869	9.63%	1.145%
17	18.839	2709	2712	2716	rBV3	47759	59471	1.12%	0.134%
18	18.986	2731	2737	2740	rBV2	65947	115771	2.19%	0.260%
19	19.121	2755	2760	2770	rBV4	215505	398978	7.53%	0.896%
20	19.345	2794	2798	2802	rBV	52163	75511	1.43%	0.170%
21	19.386	2802	2805	2810	rBV	60901	77535	1.46%	0.174%
22	19.562	2829	2835	2842	rBV	4526826	5296940	100.00%	11.897%
23	19.697	2854	2858	2862	rVB3	109966	131196	2.48%	0.295%
24	19.744	2863	2866	2873	rVB9	41699	74241	1.40%	0.167%
25	19.839	2878	2882	2887	rBV	276351	362990	6.85%	0.815%
26	20.044	2914	2917	2922	rVB7	43301	62933	1.19%	0.141%
27	20.127	2926	2931	2934	rVV	314004	386352	7.29%	0.868%
28	20.174	2936	2939	2943	rBV6	71727	88564	1.67%	0.199%
29	20.280	2952	2957	2962	rVB4	151606	241435	4.56%	0.542%
30	20.380	2970	2974	2977	rBV5	59391	78294	1.48%	0.176%
31	20.444	2978	2985	2990	rVV2	192544	368039	6.95%	0.827%
32	20.592	3006	3010	3016	rBV2	182725	222933	4.21%	0.501%
33	20.650	3017	3020	3024	rBV4	86624	96471	1.82%	0.217%
34	20.850	3050	3054	3060	rVB2	324113	486417	9.18%	1.093%

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Integration Parameters: rteint.p
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

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35	20.915	3062	3065	3070	rVB6	112607	134820	2.55%	0.303%
36	21.127	3096	3101	3111	rBV	1194006	1967730	37.15%	4.420%
37	21.286	3126	3128	3132	rBV2	73643	82085	1.55%	0.184%
38	21.456	3154	3157	3162	rBV6	158450	203847	3.85%	0.458%
39	21.709	3198	3200	3202	rBV	88043	83651	1.58%	0.188%
40	22.156	3272	3276	3282	rVB2	342288	410668	7.75%	0.922%
41	22.633	3354	3357	3362	rBV2	163234	245361	4.63%	0.551%
42	23.285	3462	3468	3481	rVB	855115	1612352	30.44%	3.621%

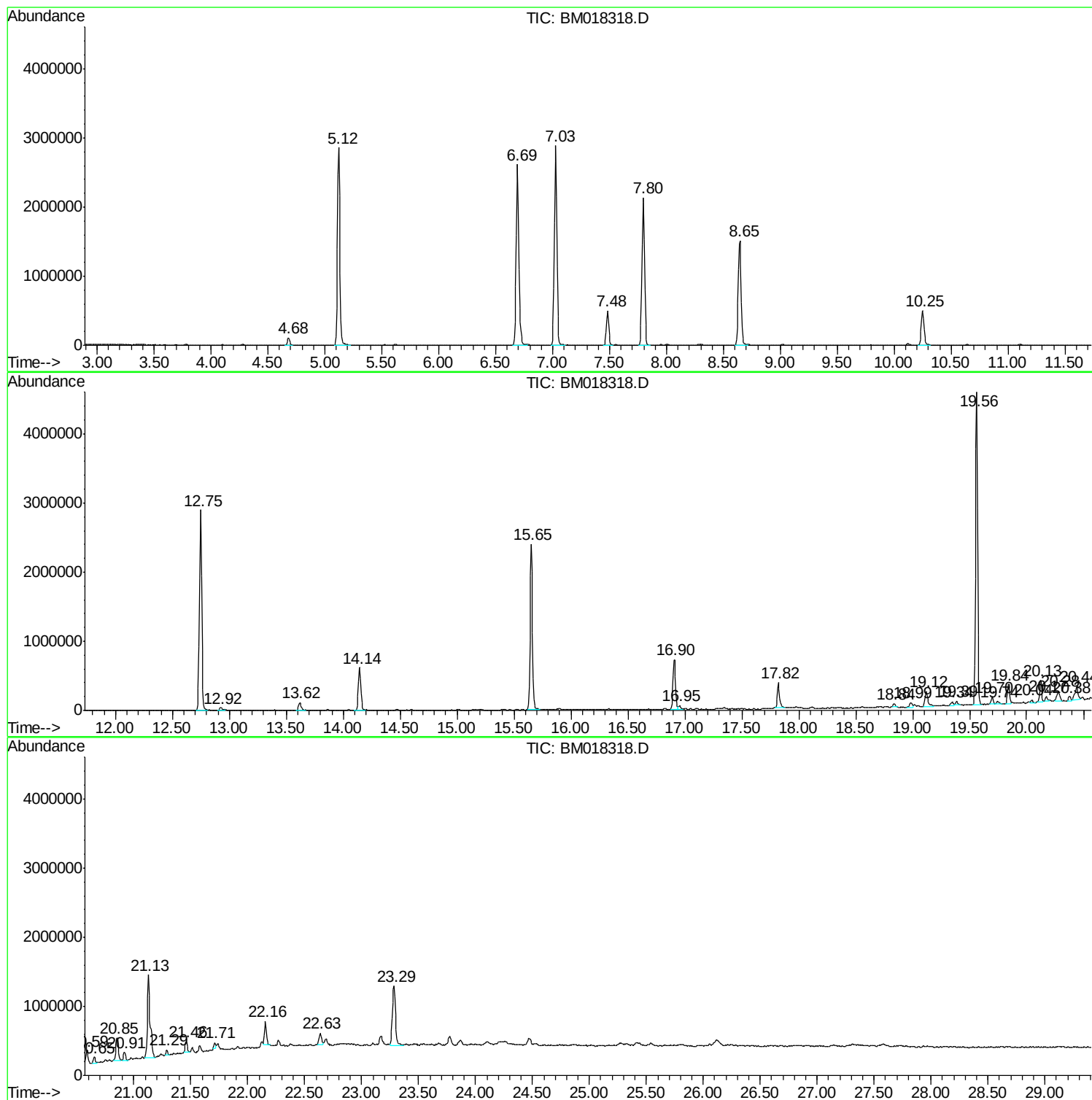
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM121518.M
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



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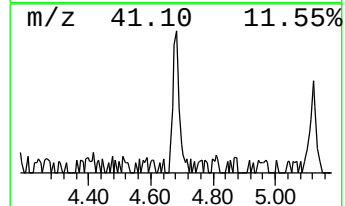
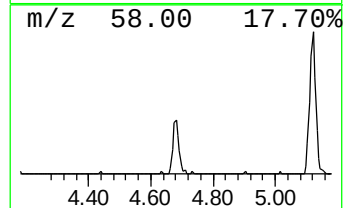
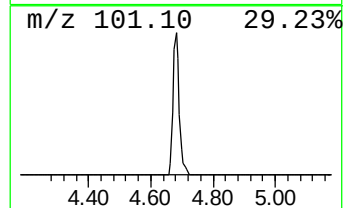
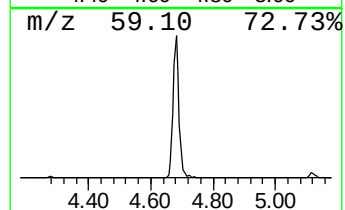
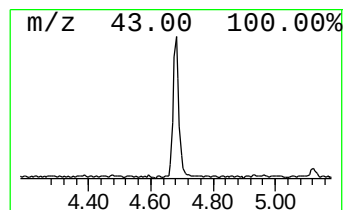
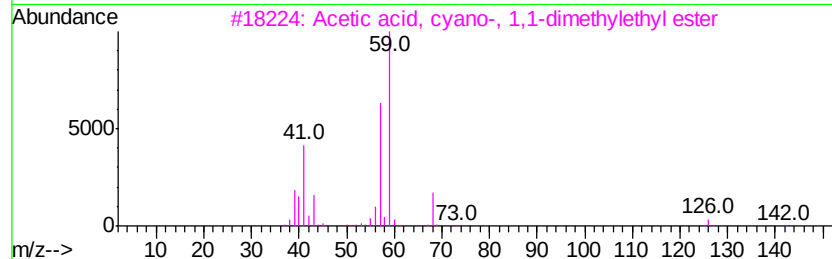
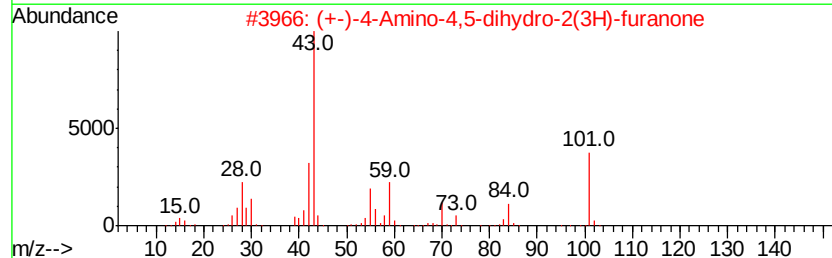
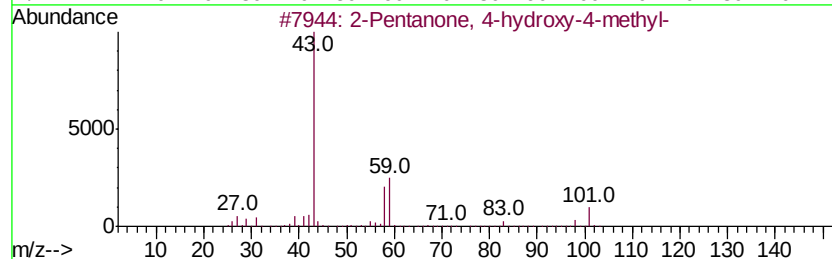
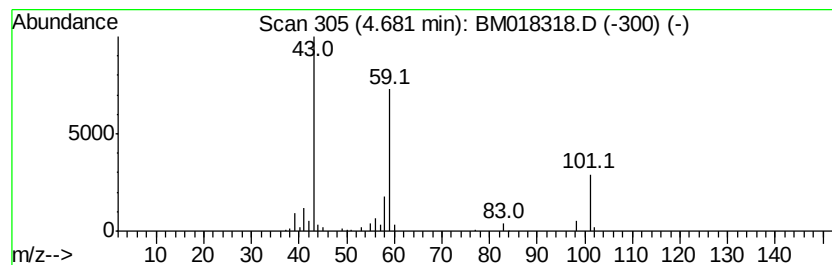
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.68	3.82 ng	146473	1,4-Dichlorobenzene-d4	7.48

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2			(+)-4-Amino-4,5-dihydro-2(3H)-f...	101	C4H7NO2	016504-58-8	47
3			Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	17
4			t-Butyl ethyl ether	102	C6H14O	1000118-18-4	12
5			3-Hydroxy-3-methyl-2-butanone	102	C5H10O2	000115-22-0	12



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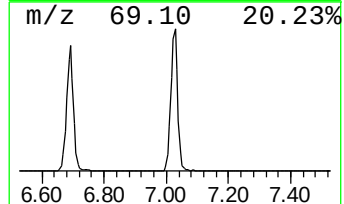
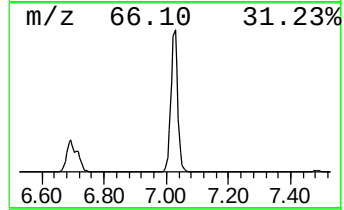
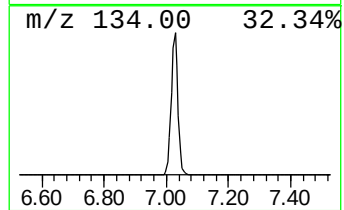
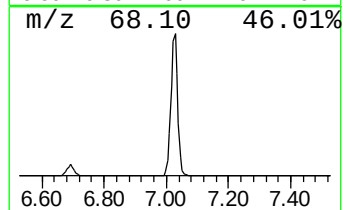
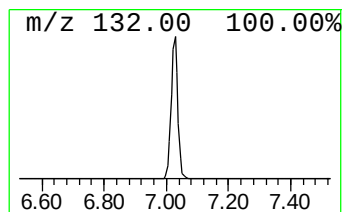
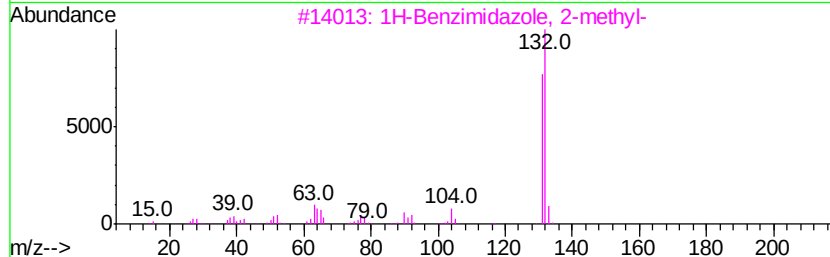
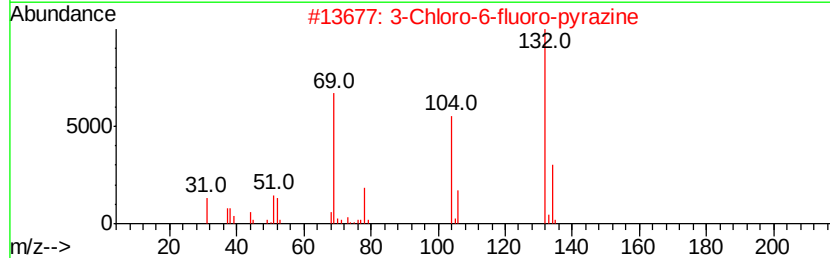
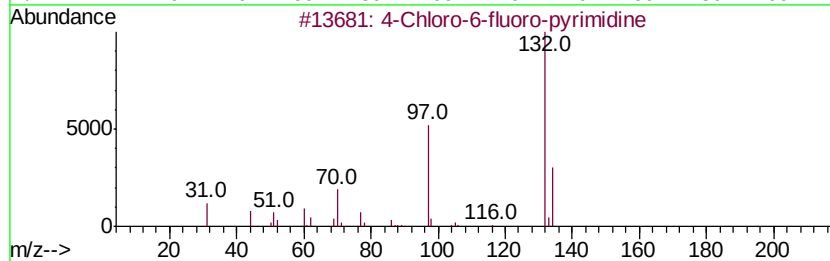
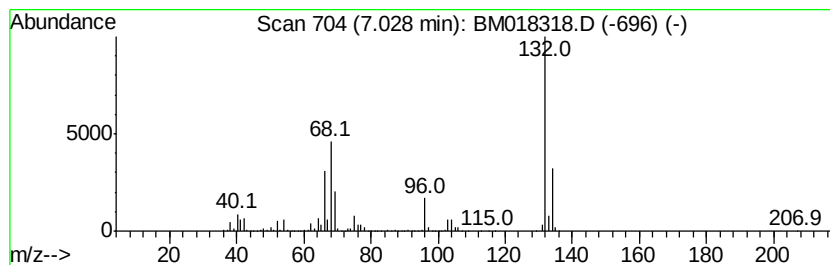
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Peak Number 2 unknown7.03 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.03	114.55 ng	4390800	1,4-Dichlorobenzene-d4	7.48

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	27
2		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
3		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
4		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
5		(5-METHYL-2-PYRIDYL)ACETONITRILE	132	C8H8N2	1000241-93-9	10



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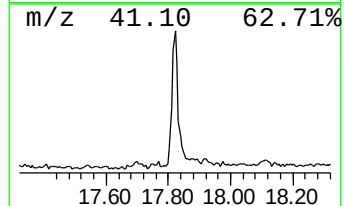
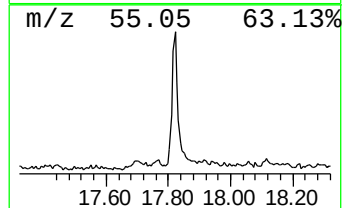
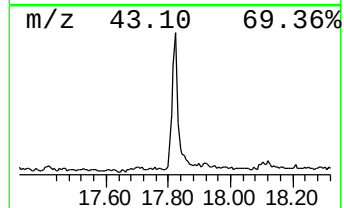
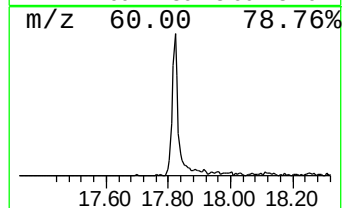
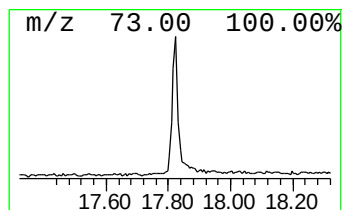
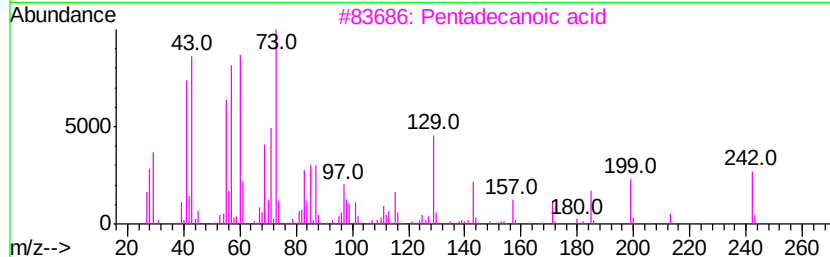
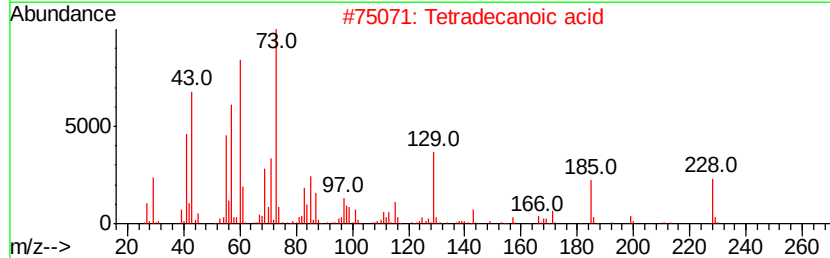
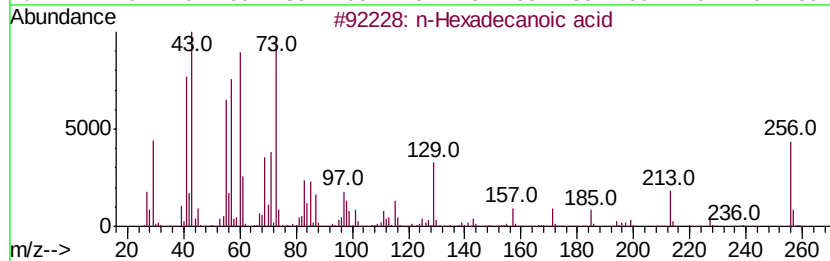
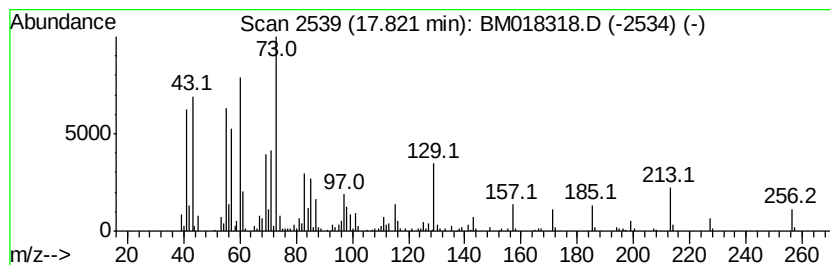
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.	
17.82	9.30 ng	509869	Phenanthrene-d10	16.91	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
2	Tetradecanoic acid	228	C14H28O2	000544-63-8	95
3	Pentadecanoic acid	242	C15H30O2	001002-84-2	80
4	Tridecanoic acid	214	C13H26O2	000638-53-9	64
5	N-Cyclooct-4-enylacetamide	167	C10H17NO	170952-69-9	25



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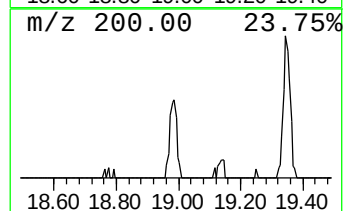
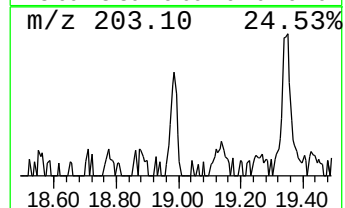
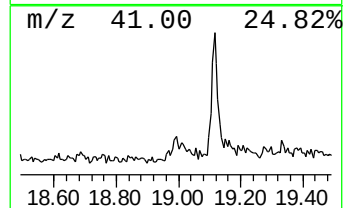
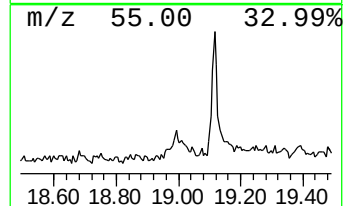
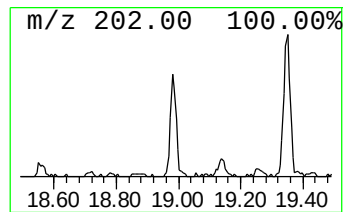
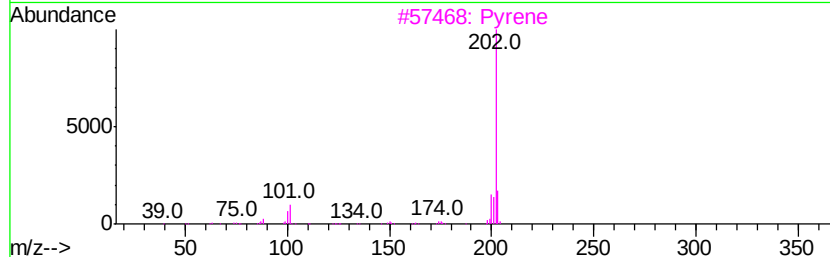
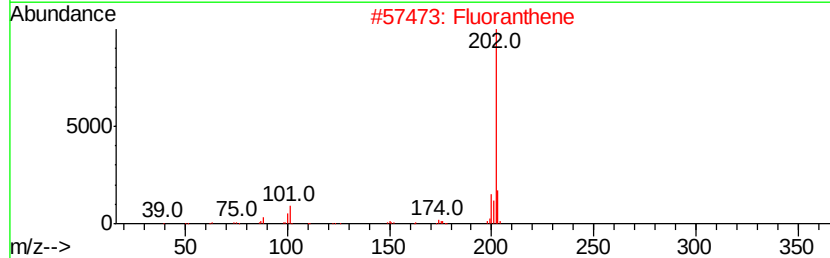
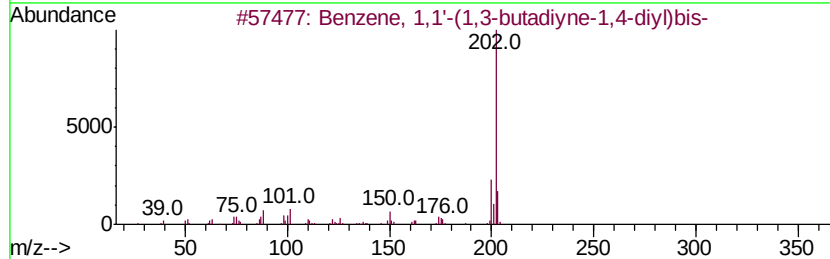
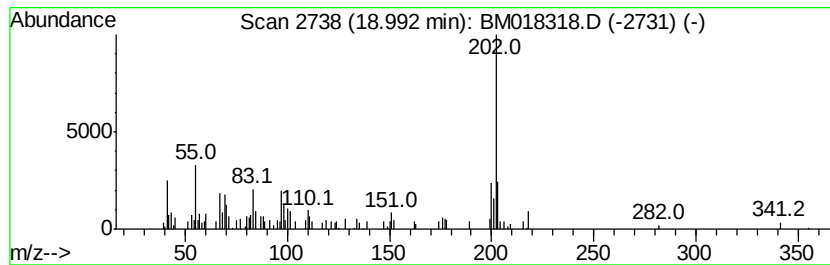
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TIC Integration Parameters: LSCINT.P

Peak Number 5 Benzene, 1,1'-(1,3-butadiyn... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.	
18.99	2.11 ng	115771	Phenanthrene-d10	16.91	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,1'-(1,3-butadiyne-1,4...	202	C16H10	000886-66-8	86
2	Fluoranthene	202	C16H10	000206-44-0	58
3	Pvrene	202	C16H10	000129-00-0	58
4	2H-Isoindole-2-acetic acid, 1,3-...	247	C13H13N04	019506-85-5	47
5	1,4-Pentadiyn-3-one, 1,5-diphenyl-	230	C17H10O	015814-30-9	45



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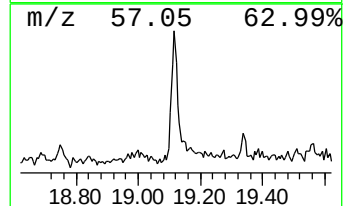
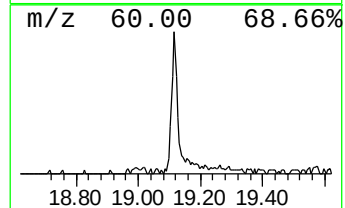
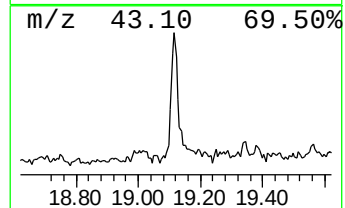
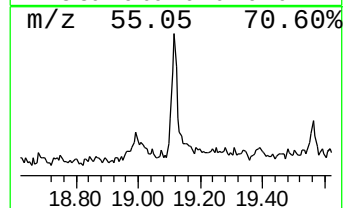
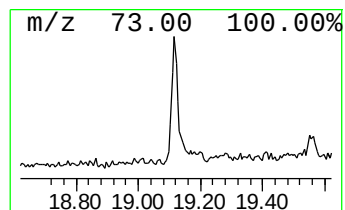
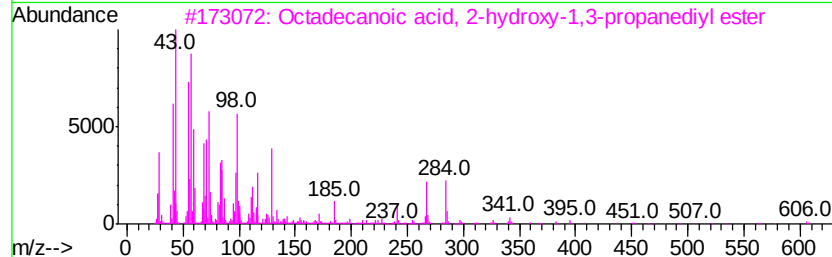
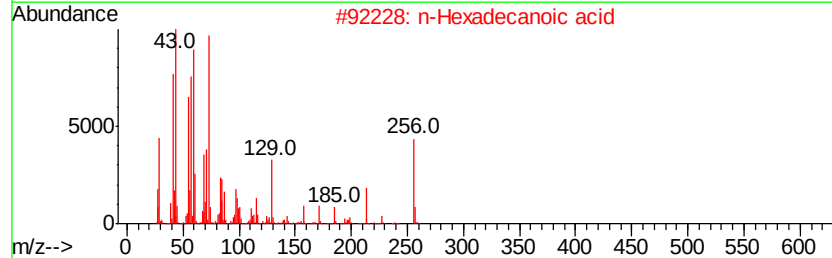
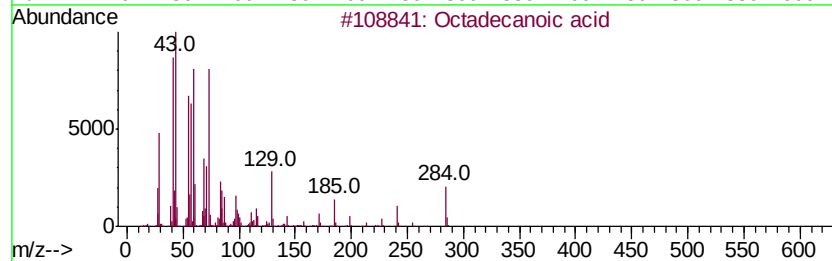
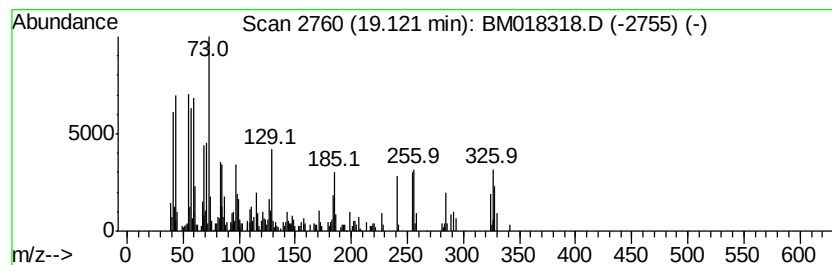
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM121518.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Octadecanoic acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
19.12	4.06 ng	398978	Chrysene-d12	21.13	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecanoic acid	284	C18H36O2	000057-11-4	89
2	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	60
3	Octadecanoic acid, 2-hydroxy-1,3...	625	C39H76O5	000504-40-5	35
4	Tetradecanoic acid	228	C14H28O2	000544-63-8	30
5	Pentadecanoic acid	242	C15H30O2	001002-84-2	30



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM121918\
Data File : BM018318.D
Acq On : 20 Dec 2018 04:14
Operator : JU/SJ
Sample : J6378-24
Misc :
ALS Vial : 19 Sample Multiplier: 1

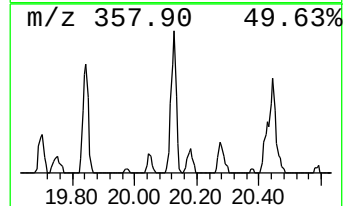
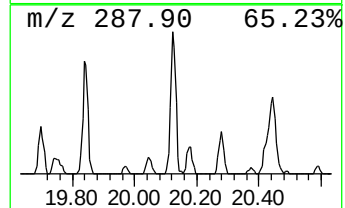
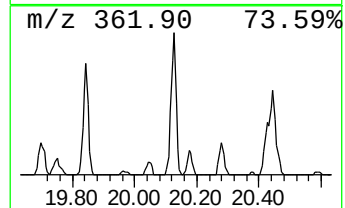
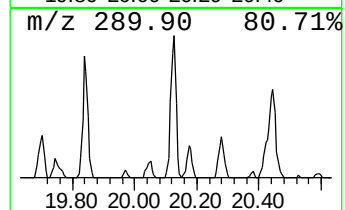
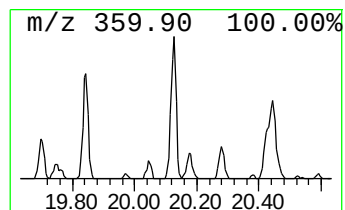
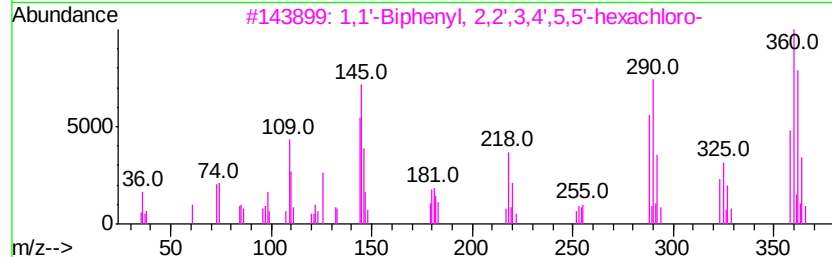
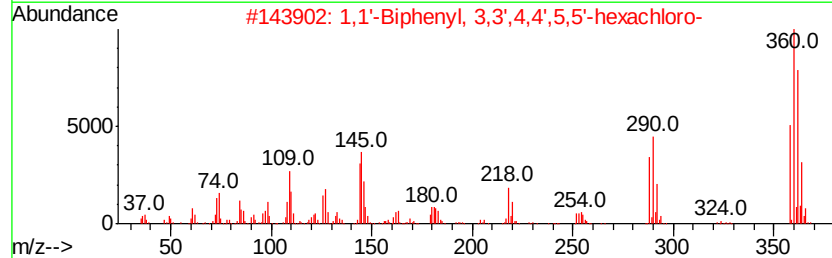
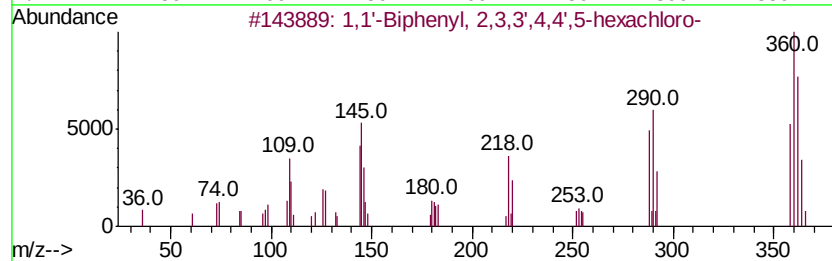
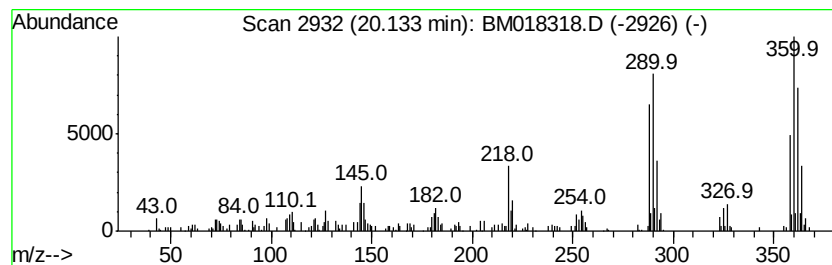
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 1,1'-Biphenyl, 2,3,3',4,4',... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.	
20.13	3.93 ng	386352	Chrysene-d12	21.13	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,3,3',4,4',5-hex...	358	C12H4Cl6	038380-08-4	98
2	1,1'-Biphenyl, 3,3',4,4',5,5'-he...	358	C12H4Cl6	032774-16-6	98
3	1,1'-Biphenyl, 2,2',3,4',5,5'-he...	358	C12H4Cl6	051908-16-8	96
4	1,1'-Biphenyl, 2,2',3,3',4,5-hex...	358	C12H4Cl6	055215-18-4	96
5	1,1'-Biphenyl, 2,2',3,3',6,6'-he...	358	C12H4Cl6	038411-22-2	96



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM121918\
Data File : BM018318.D
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ClientSampleId :
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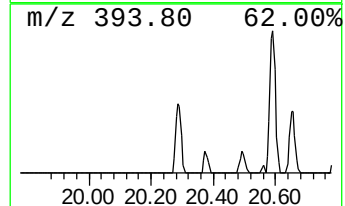
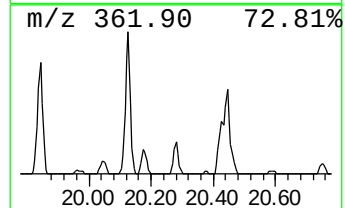
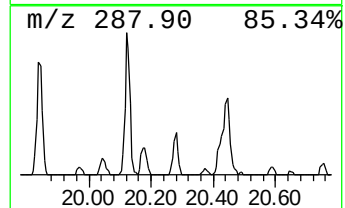
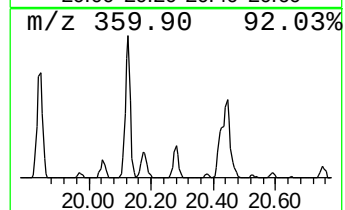
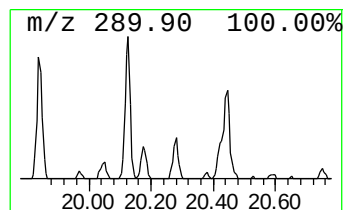
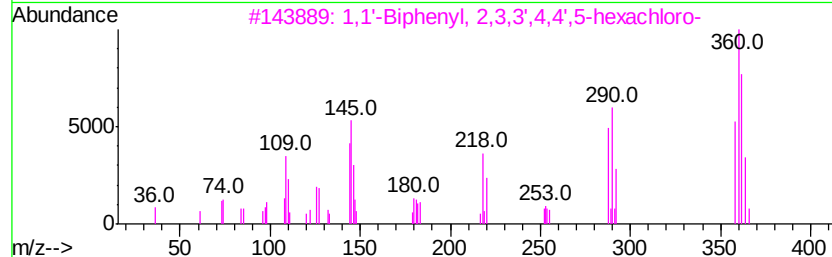
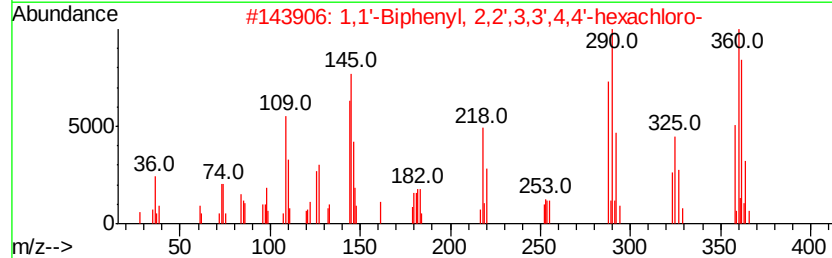
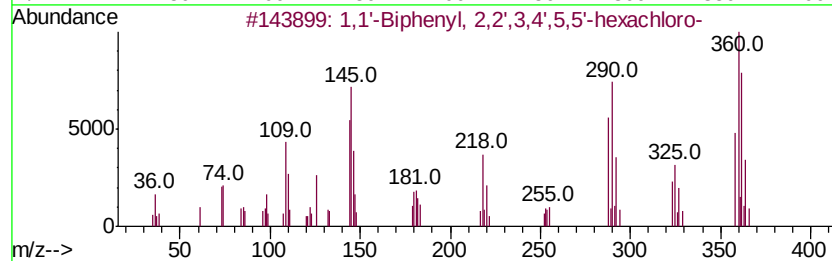
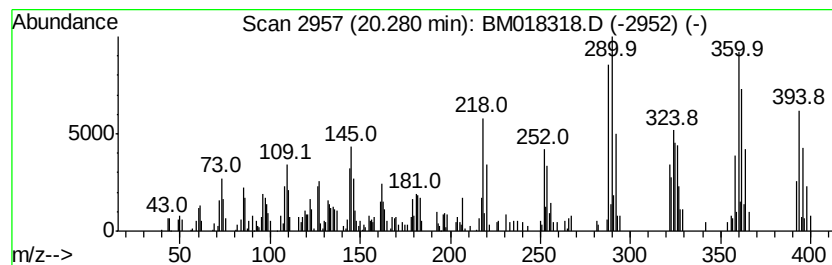
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 1,1'-Biphenyl, 2,2',3,4',5,... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.28	2.45 ng	241435	Chrysene-d12	21.13

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,1'-Biphenyl, 2,2',3,4',5,5'-he...	358	C12H4Cl6	051908-16-8	99
2		1,1'-Biphenyl, 2,2',3,3',4,4'-he...	358	C12H4Cl6	038380-07-3	99
3		1,1'-Biphenyl, 2,3,3',4,4',5-hex...	358	C12H4Cl6	038380-08-4	99
4		1,1'-Biphenyl, 2,2',3,3',4,5'-he...	358	C12H4Cl6	052663-66-8	98
5		1,1'-Biphenyl, 2,2',3,3',4,5-hex...	358	C12H4Cl6	055215-18-4	97



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM121918\
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Operator : JU/SJ
Sample : J6378-24
Misc :
ALS Vial : 19 Sample Multiplier: 1

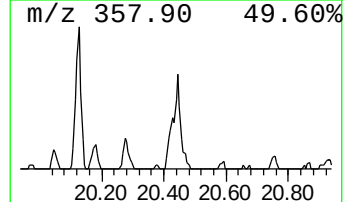
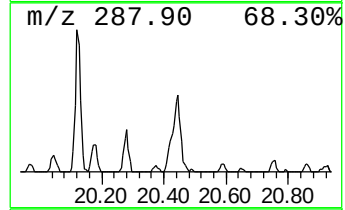
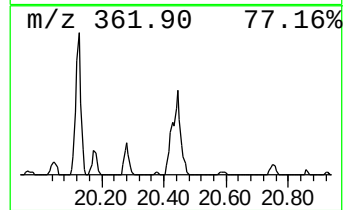
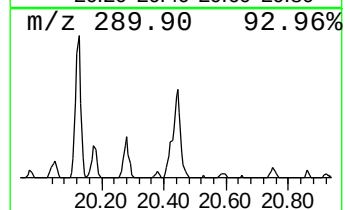
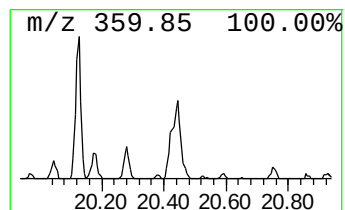
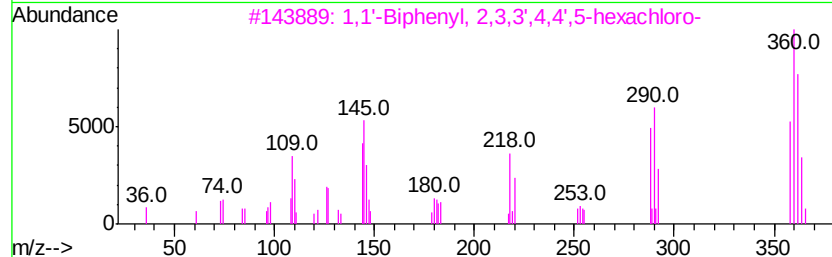
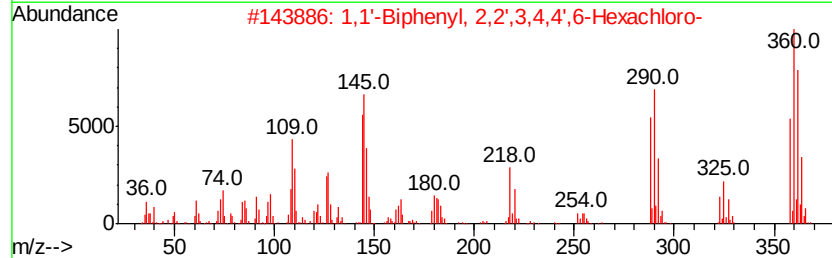
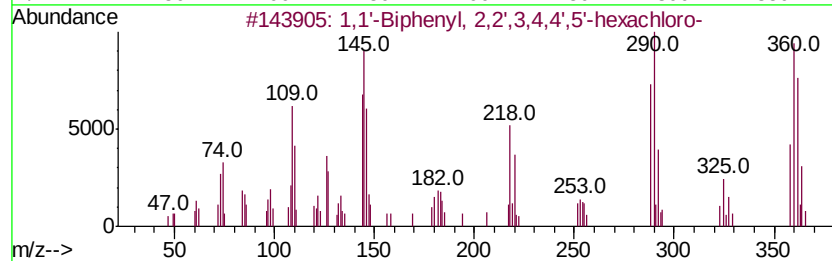
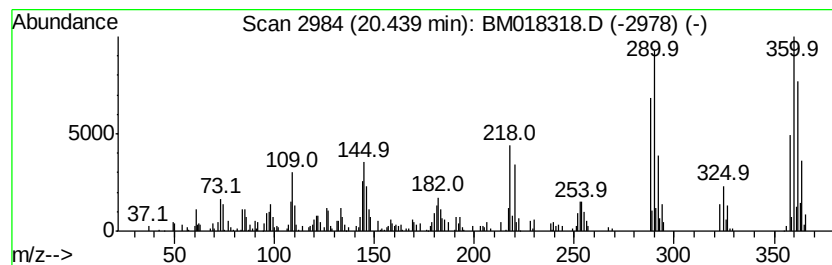
Instrument :
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM121518.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 1,1'-Biphenyl, 2,2',3,4,4',... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.		
20.44	3.74 ng	368039	Chrysene-d12	21.13		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl,	2,2',3,4,4',5'-he...	358	C12H4Cl6	035065-28-2	99
2	1,1'-Biphenyl,	2,2',3,4,4',6-Hex...	358	C12H4Cl6	056030-56-9	97
3	1,1'-Biphenyl,	2,3,3',4,4',5-hex...	358	C12H4Cl6	038380-08-4	96
4	1,1'-Biphenyl,	2,2',3,3',5,6'-he...	358	C12H4Cl6	052744-13-5	96
5	1,1'-Biphenyl,	2,3',4,4',5,5'-he...	358	C12H4Cl6	052663-72-6	95



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM121918\
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Acq On : 20 Dec 2018 04:14
Operator : JU/SJ
Sample : J6378-24
Misc :
ALS Vial : 19 Sample Multiplier: 1

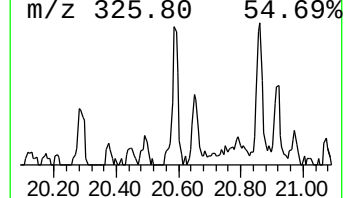
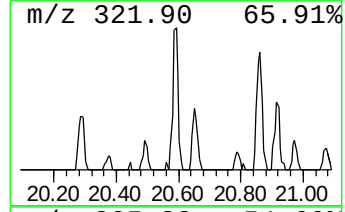
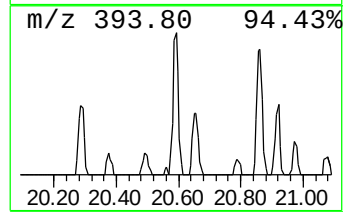
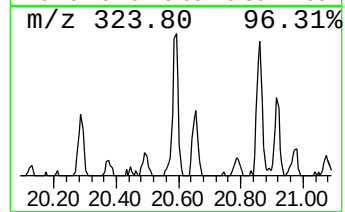
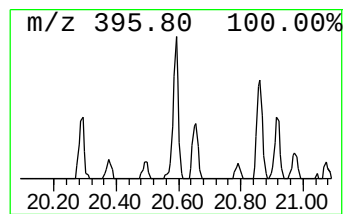
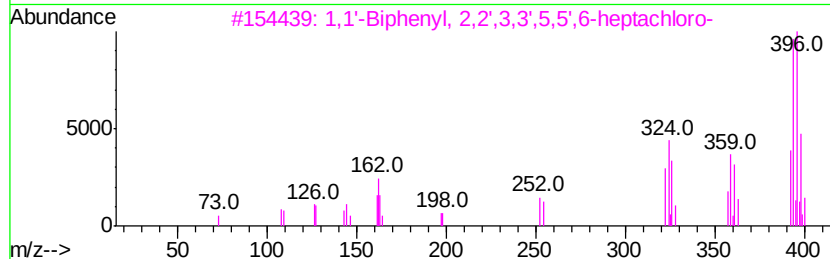
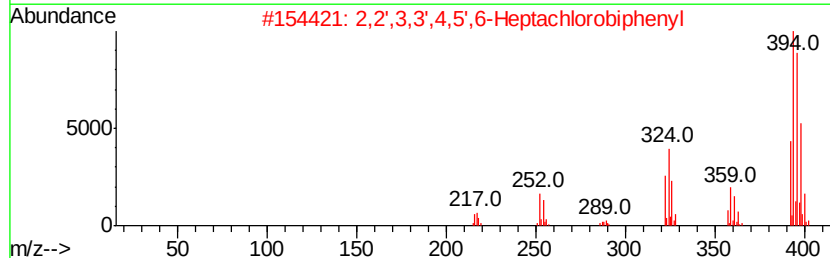
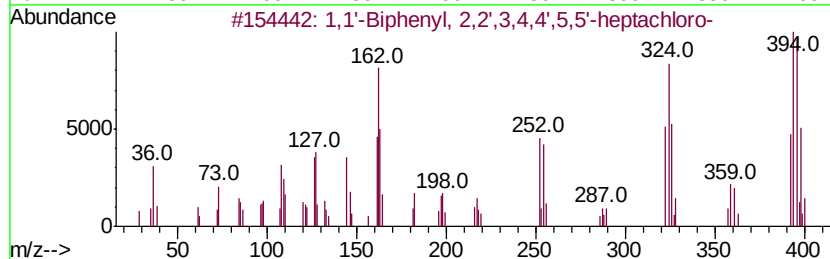
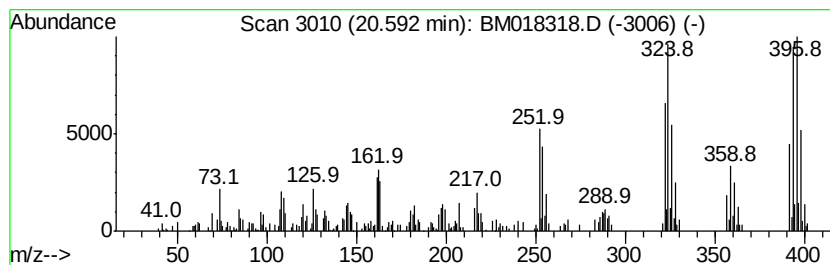
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM121518.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 1,1'-Biphenyl, 2,2',3,4,4',... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.	
20.59	2.27 ng	222933	Chrysene-d12	21.13	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,2',3,4,4',5,5'-...	392	C12H3Cl7	035065-29-3	99
2	2,2',3,3',4,5',6-Heptachlorobiph...	392	C12H3Cl7	040186-70-7	99
3	1,1'-Biphenyl, 2,2',3,3',5,5',6-...	392	C12H3Cl7	052663-67-9	99
4	1,1'-Biphenyl, 2,2',3,4,5,6,6'-H...	392	C12H3Cl7	074472-49-4	99
5	1,1'-Biphenyl, 2,3,3',4,5,5',6-h...	392	C12H3Cl7	074472-51-8	99



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM121918\
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Acq On : 20 Dec 2018 04:14
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Sample : J6378-24
Misc :
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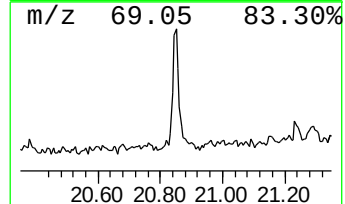
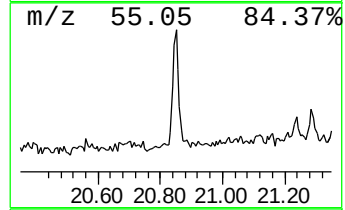
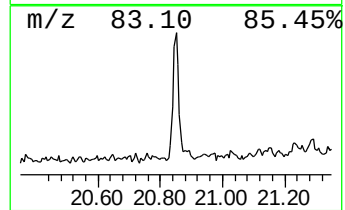
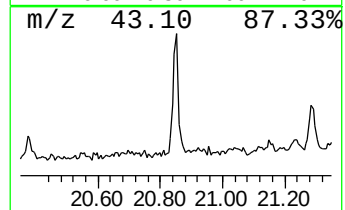
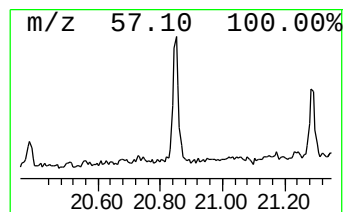
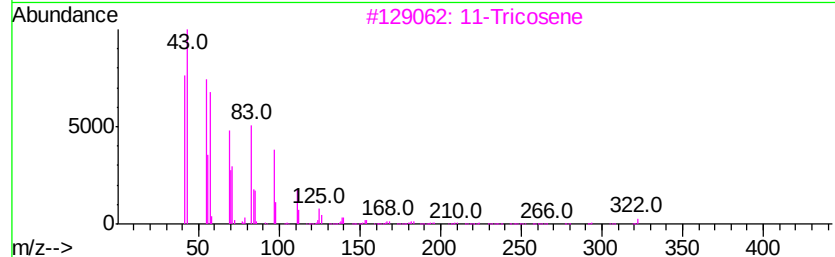
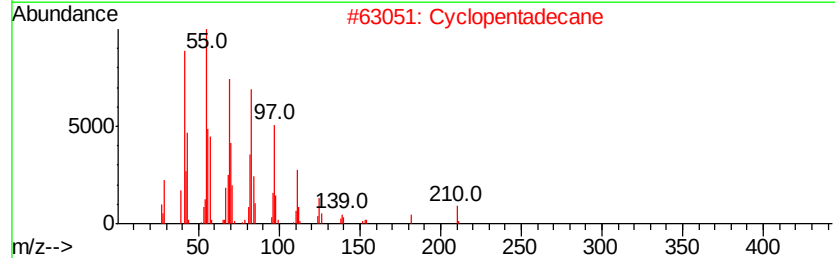
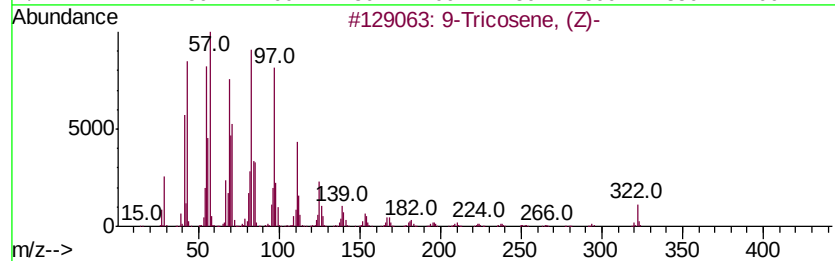
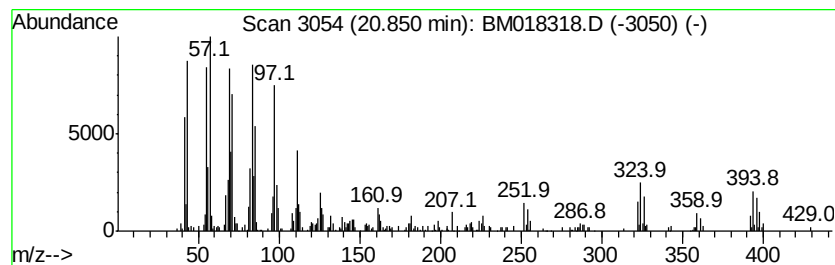
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM121518.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 12 9-Tricosene, (Z)- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.	
20.85	4.94 ng	486417	Chrysene-d12	21.13	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9-Tricosene, (Z)-	322	C23H46	027519-02-4	97
2	Cyclopentadecane	210	C15H30	000295-48-7	97
3	11-Tricosene	322	C23H46	052078-56-5	93
4	1-Tricosene	322	C23H46	018835-32-0	91
5	Cyclooctacosane	392	C28H56	000297-24-5	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM121918\
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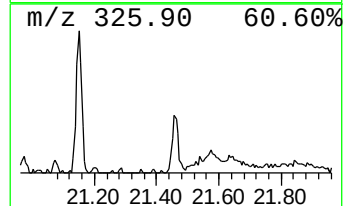
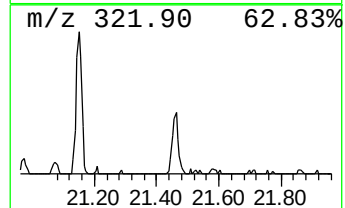
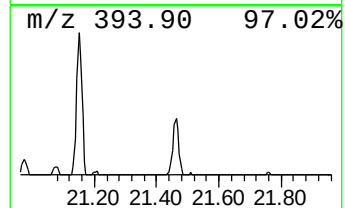
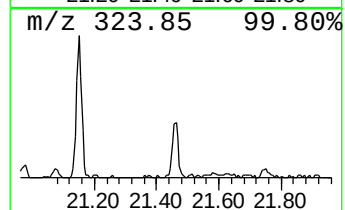
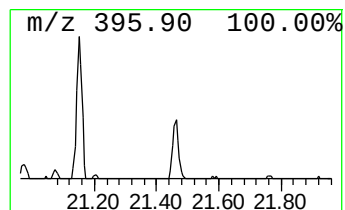
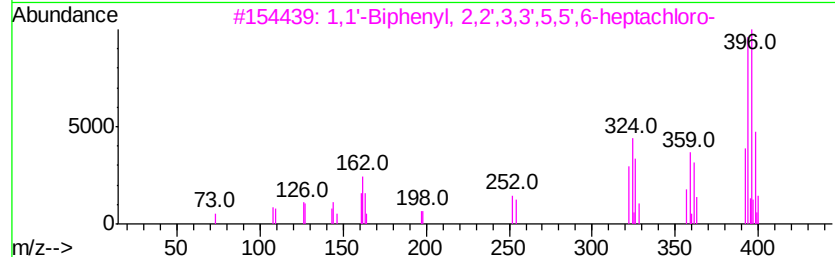
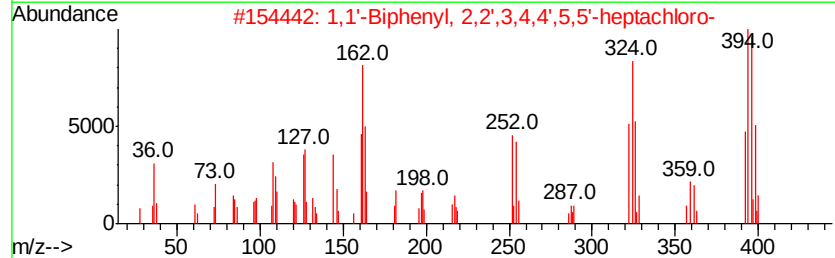
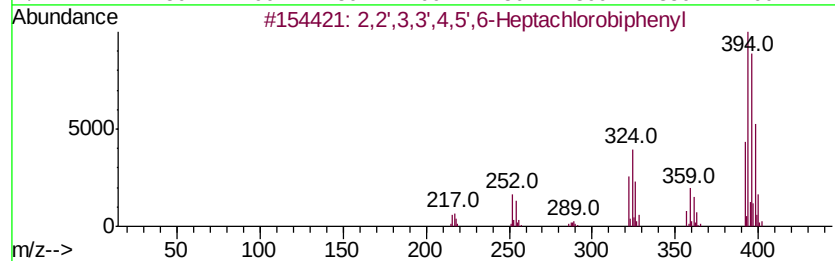
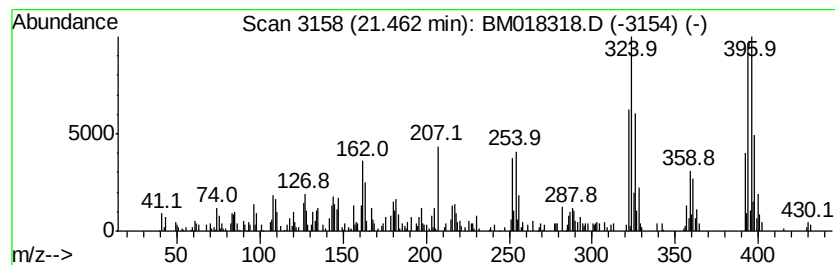
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 13 2,2',3,3',4,5',6-Heptachlor... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.		
21.46	2.07 ng	203847	Chrysene-d12	21.13		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,2',3,3',4,5',6-Heptachlorobiph...	392	C12H3Cl7	040186-70-7	99	
2	1,1'-Biphenyl, 2,2',3,4,4',5,5'-...	392	C12H3Cl7	035065-29-3	99	
3	1,1'-Biphenyl, 2,2',3,3',5,5',6-...	392	C12H3Cl7	052663-67-9	99	
4	1,1'-Biphenyl, 2,3,3',4,5,5',6-h...	392	C12H3Cl7	074472-51-8	99	
5	1,1'-Biphenyl, 2,2',3,4,4',5,6-H...	392	C12H3Cl7	074472-47-2	99	



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM121918\
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Misc :
ALS Vial : 19 Sample Multiplier: 1

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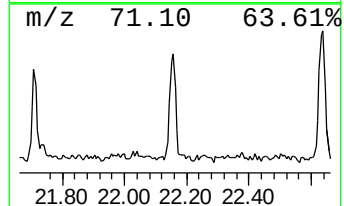
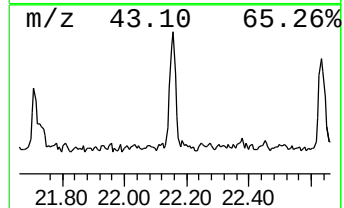
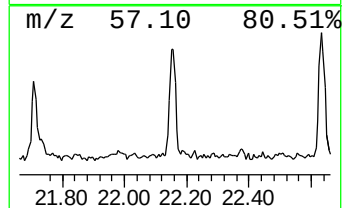
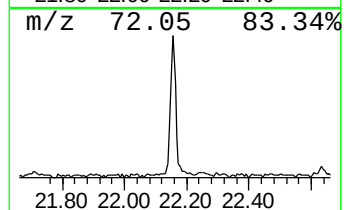
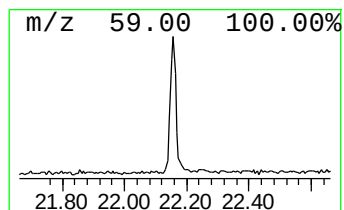
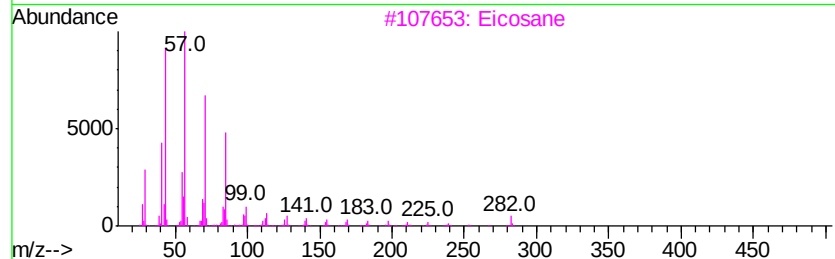
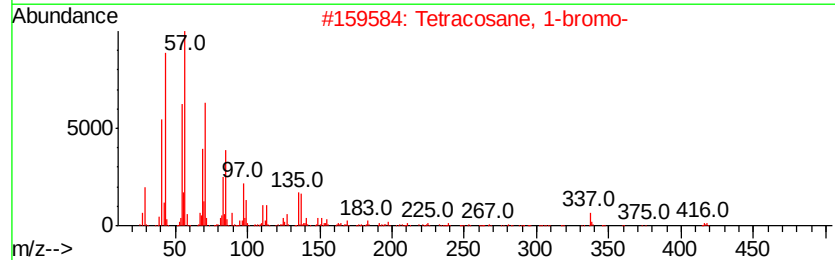
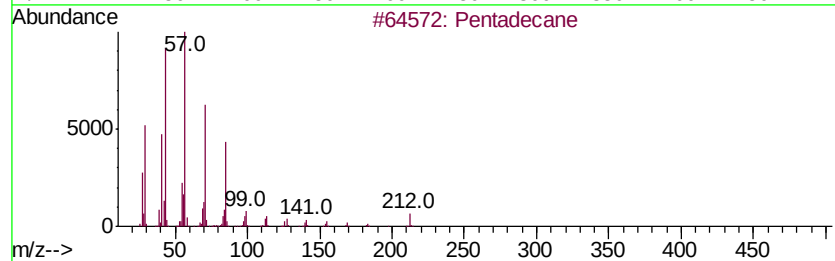
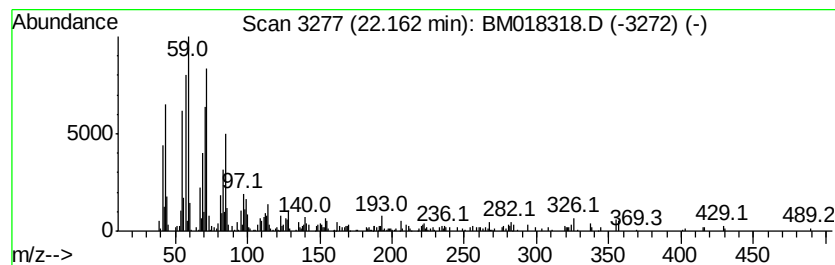
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 14 Pentadecane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.16	4.17 ng	410668	Chrysene-d12	21.13

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentadecane	212	C15H32	000629-62-9	55
2		Tetracosane, 1-bromo-	416	C24H49Br	006946-24-3	51
3		Eicosane	282	C20H42	000112-95-8	35
4		Nonadecane	268	C19H40	000629-92-5	35
5		Octadecane	254	C18H38	000593-45-3	35



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM121918\
Data File : BM018318.D
Acq On : 20 Dec 2018 04:14
Operator : JU/SJ
Sample : J6378-24
Misc :
ALS Vial : 19 Sample Multiplier: 1

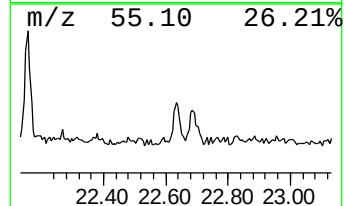
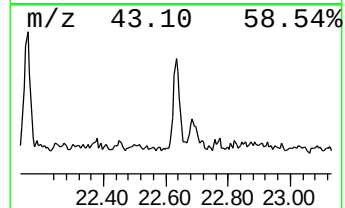
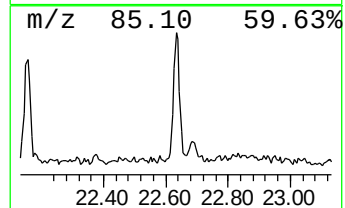
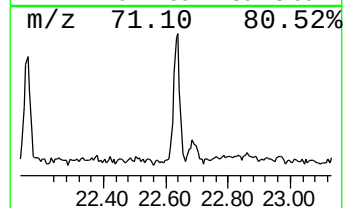
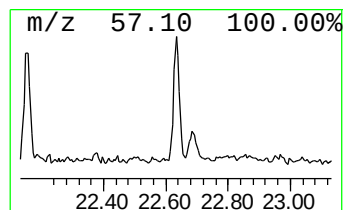
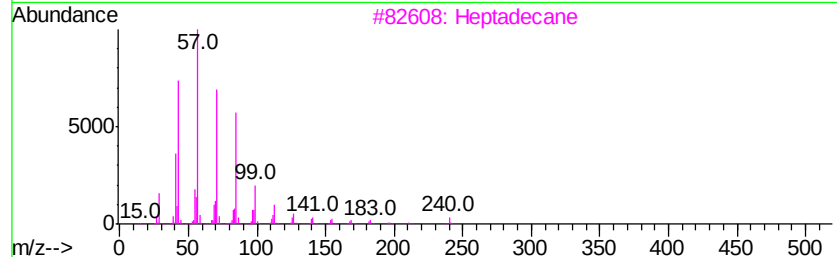
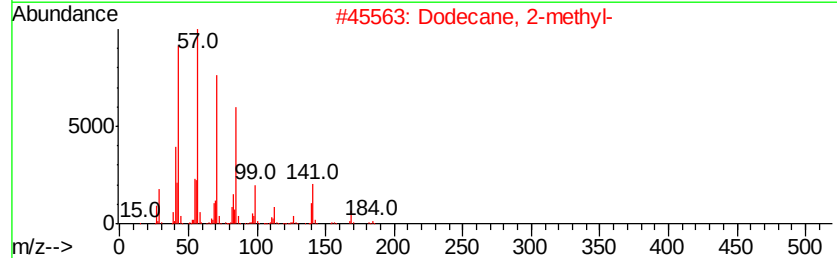
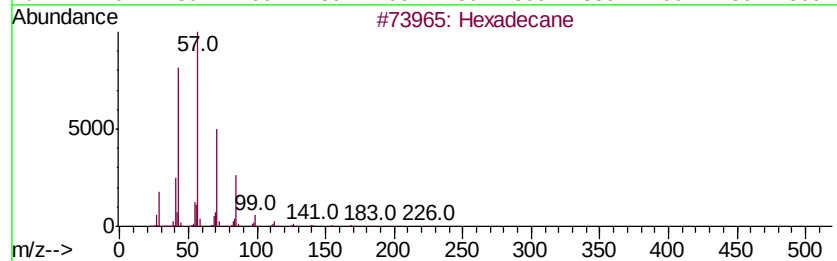
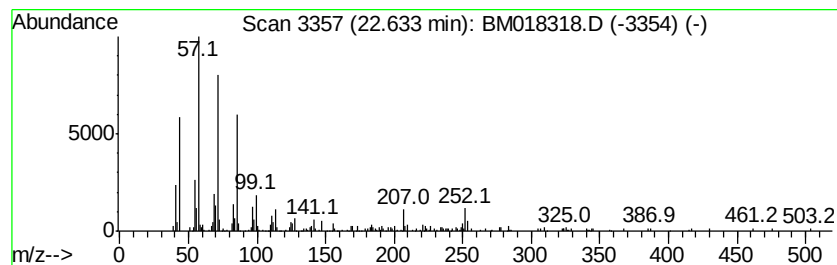
Instrument :
BNA_M
ClientSampleId :
P001-SS002-1218-01

Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM121518.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 15 Hexadecane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD		R.T.
22.63	3.04 ng	245361	Perylene-d12		23.29
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane	226	C16H34	000544-76-3	94
2	Dodecane, 2-methyl-	184	C13H28	001560-97-0	76
3	Heptadecane	240	C17H36	000629-78-7	76
4	Heneicosane	296	C21H44	000629-94-7	76
5	Octadecane	254	C18H38	000593-45-3	68



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM121918\
 Data File : BM018318.D
 Acq On : 20 Dec 2018 04:14
 Operator : JU/SJ
 Sample : J6378-24
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 P001-SS002-1218-01

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\8270-BM121518.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
 TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	---Internal Standard---			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.68	3.8	ng	146473	1	7.48	766642	20.0
unknown7.03	7.03	114.6	ng	4390800	1	7.48	766642	20.0
n-Hexadecanoic acid	17.82	9.3	ng	509869	4	16.91	1096020	20.0
Benzene, 1,1'-(1,...	18.99	2.1	ng	115771	4	16.91	1096020	20.0
Octadecanoic acid	19.12	4.1	ng	398978	5	21.13	1967730	20.0
1,1'-Biphenyl, 2,...	20.13	3.9	ng	386352	5	21.13	1967730	20.0
1,1'-Biphenyl, 2,...	20.28	2.5	ng	241435	5	21.13	1967730	20.0
1,1'-Biphenyl, 2,...	20.44	3.7	ng	368039	5	21.13	1967730	20.0
1,1'-Biphenyl, 2,...	20.59	2.3	ng	222933	5	21.13	1967730	20.0
9-Tricosene, (Z)-	20.85	4.9	ng	486417	5	21.13	1967730	20.0
2,2',3,3',4,5',6-...	21.46	2.1	ng	203847	5	21.13	1967730	20.0
Pentadecane	22.16	4.2	ng	410668	5	21.13	1967730	20.0
Hexadecane	22.63	3.0	ng	245361	6	23.29	1612350	20.0