

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG012020\
 Data File : BG044097.D
 Acq On : 20 Jan 2020 14:27
 Operator : JU
 Sample : L1176-03
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 EP-3

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_G\METHODS\8270-BG123019.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	5.034	325	331	342	rBV	222973	342597	8.74%	1.258%
2	5.510	405	412	431	rBV	978738	1606905	41.01%	5.899%
3	7.102	674	683	694	rBV	1076342	1958881	49.99%	7.192%
4	7.467	737	745	756	rBV	1061169	1825901	46.60%	6.703%
5	7.931	818	824	833	rBV	291690	499130	12.74%	1.832%
6	8.254	871	879	889	rBV	765666	1355868	34.60%	4.978%
7	9.088	1012	1021	1036	rBV	624073	1137825	29.04%	4.177%
8	10.733	1293	1301	1313	rBV	407141	770978	19.68%	2.830%
9	13.189	1712	1719	1737	rBV	1820664	2966099	75.70%	10.889%
10	14.024	1853	1861	1874	rBV	107978	195641	4.99%	0.718%
11	14.570	1946	1954	1969	rBV	747098	1201349	30.66%	4.410%
12	16.057	2200	2207	2222	rBV	1342892	2144989	54.74%	7.875%
13	17.314	2414	2421	2432	rBV	849222	1316090	33.59%	4.832%
14	19.929	2859	2866	2880	rBV	2971473	3918436	100.00%	14.386%
15	21.239	3083	3089	3102	rBV	159216	348736	8.90%	1.280%
16	21.562	3137	3144	3155	rBV	780043	1342665	34.27%	4.929%
17	21.773	3175	3180	3186	rBV	157254	210495	5.37%	0.773%
18	22.367	3275	3281	3292	rBV	243598	404464	10.32%	1.485%
19	23.037	3388	3395	3402	rBV	292984	548171	13.99%	2.012%
20	23.818	3521	3528	3539	rVB	273687	586603	14.97%	2.154%
21	24.670	3663	3673	3680	rBV	476516	1365992	34.86%	5.015%
22	24.735	3680	3684	3696	rVB2	214909	532416	13.59%	1.955%
23	25.827	3861	3870	3883	rVB3	151487	435494	11.11%	1.599%
24	27.144	4084	4094	4103	rBV3	68458	222800	5.69%	0.818%

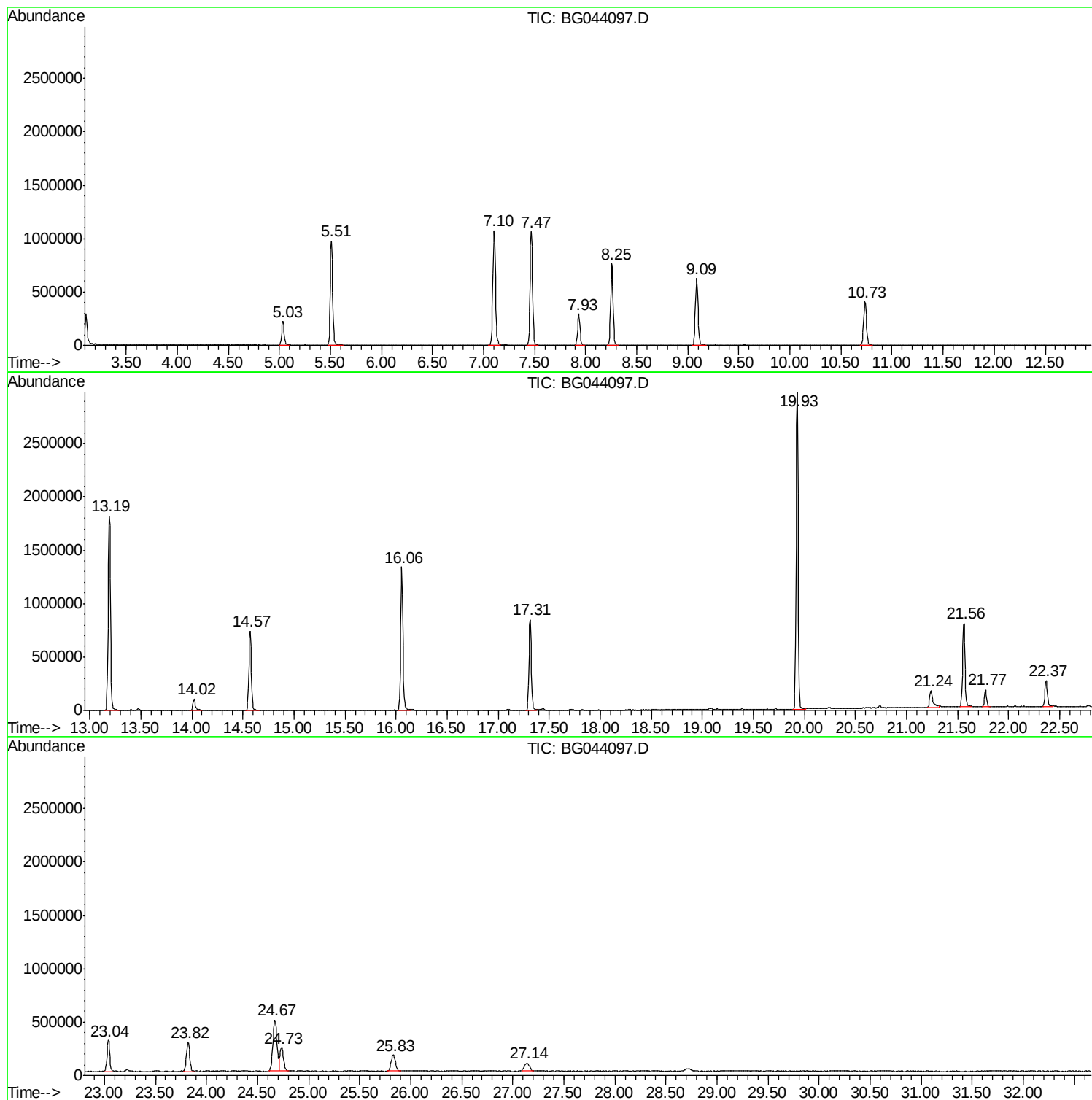
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG123019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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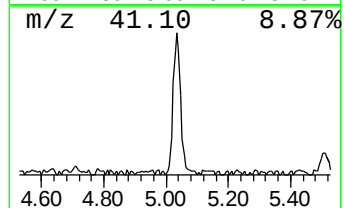
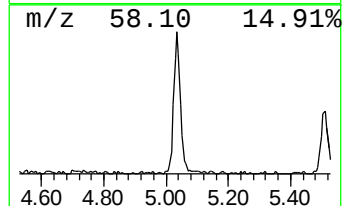
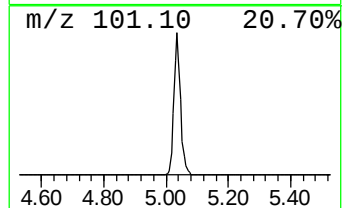
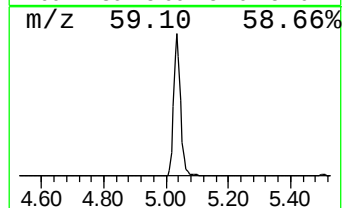
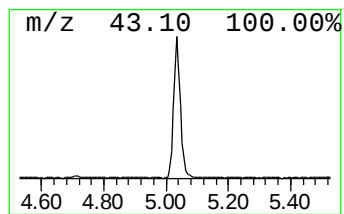
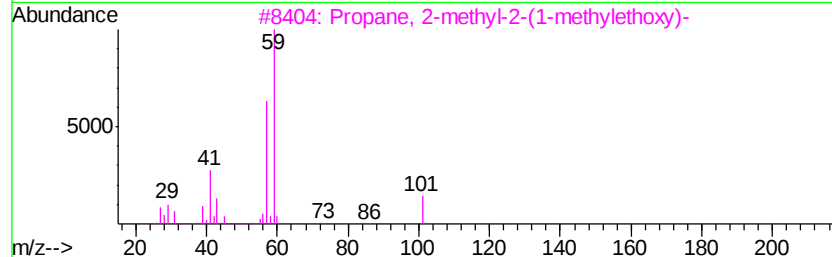
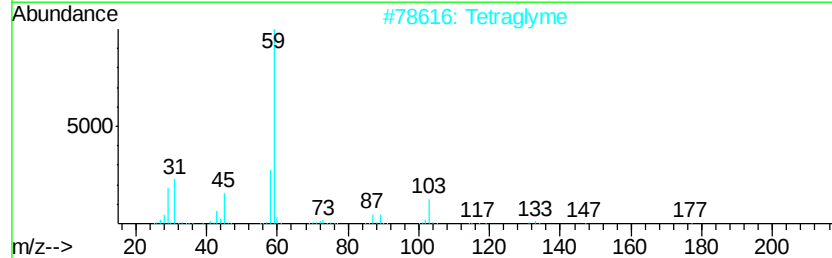
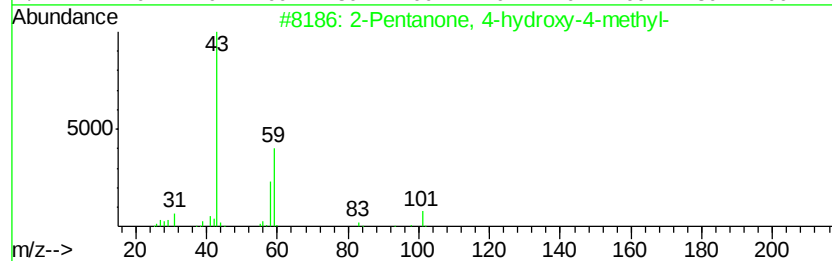
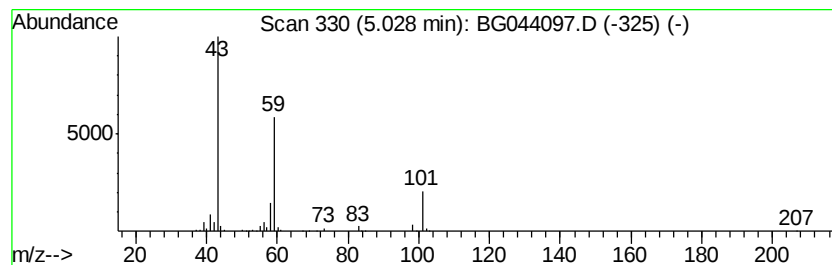
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.03	13.73 ng	342597	1,4-Dichlorobenzene-d4	7.93

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	40
2			Tetraglyme	222	C10H22O5	000143-24-8	25
3			Propane, 2-methyl-2-(1-methylethoxy)-	116	C7H16O	017348-59-3	23
4			Acetic acid, cyano-, 1,1-dimethyl-	141	C7H11NO2	001116-98-9	23
5			Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9



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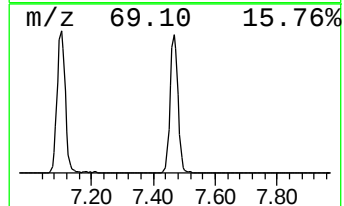
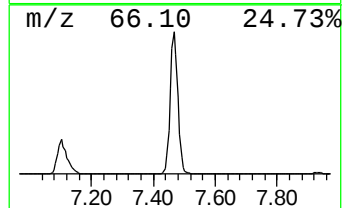
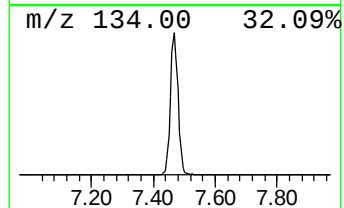
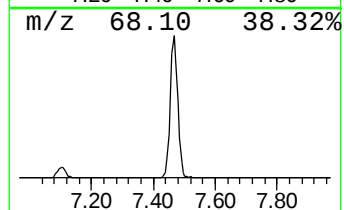
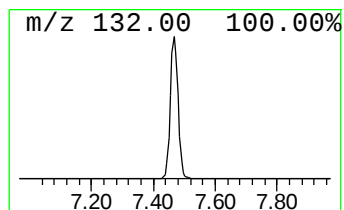
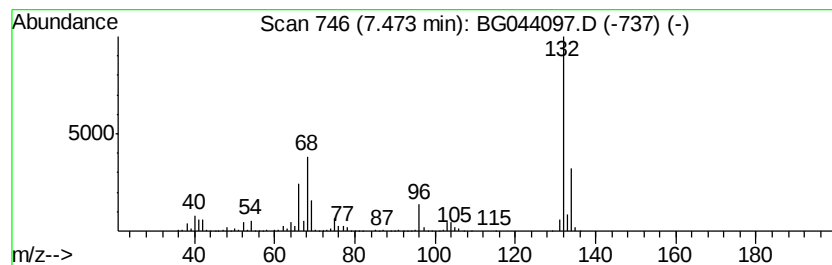
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TIC Library : C:\Database\NIST11.L
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Peak Number 2 unknown7.47 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.47	73.16 ng	1825900	1,4-Dichlorobenzene-d4	7.93

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	16
3		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	12
4		5-Aminoindole	132	C8H8N2	005192-03-0	12
5		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9



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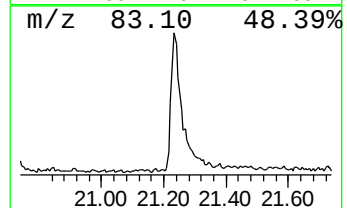
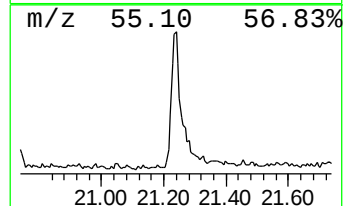
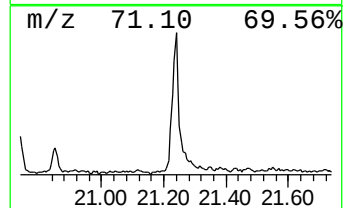
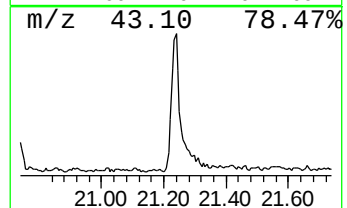
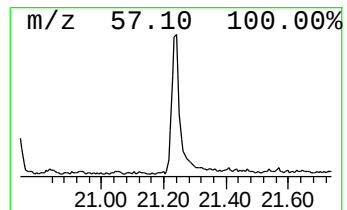
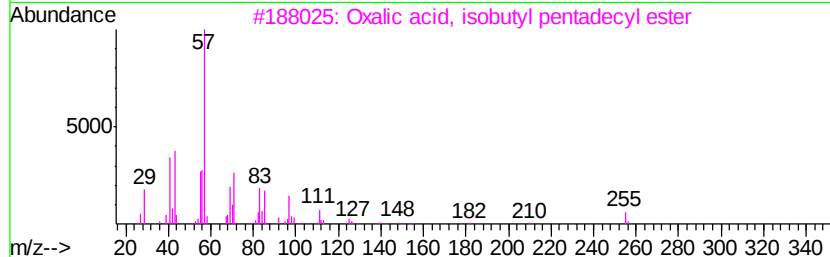
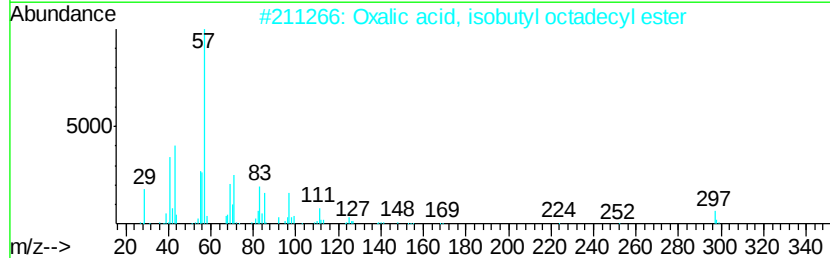
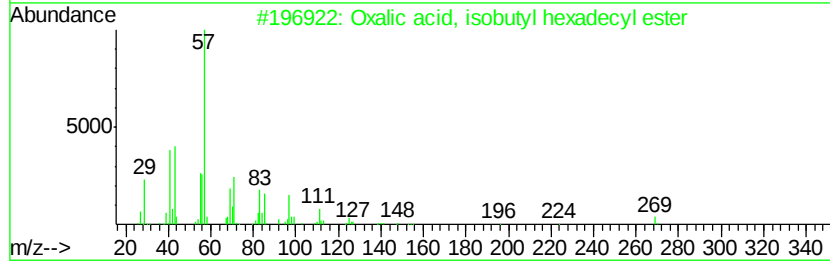
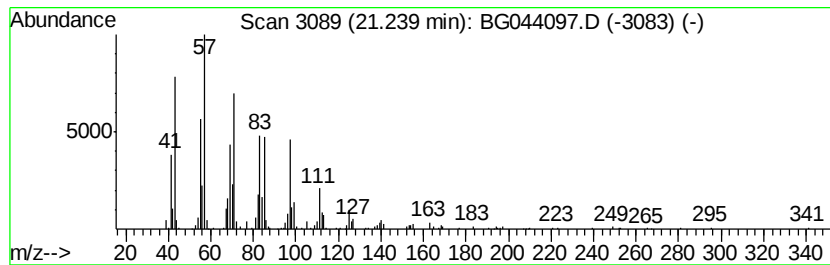
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Peak Number 4 Oxalic acid, isobutyl hexadecyl... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.24	5.19 ng	348736	Chrysene-d12	21.56

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Oxalic acid, isobutyl hexadecyl ...	370	C22H42O4	1000309-38-1	91
2			Oxalic acid, isobutyl octadecyl ...	398	C24H46O4	1000309-38-3	91
3			Oxalic acid, isobutyl pentadecyl...	356	C21H40O4	1000309-38-0	91
4			1-Decanol, 2-octyl-	270	C18H38O	045235-48-1	91
5			Heneicosyl heptafluorobutyrate	508	C25H43F7O2	1000351-83-8	81



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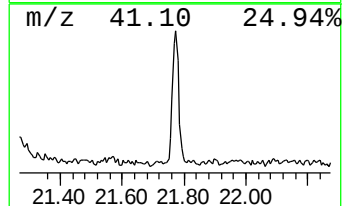
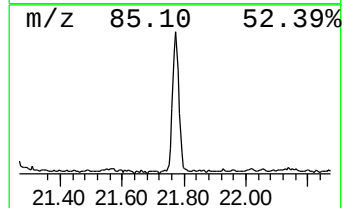
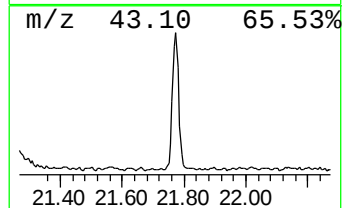
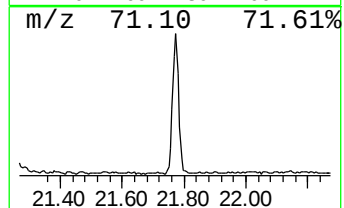
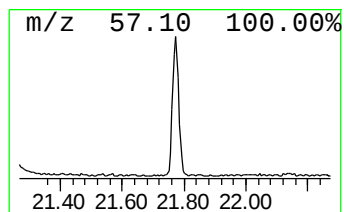
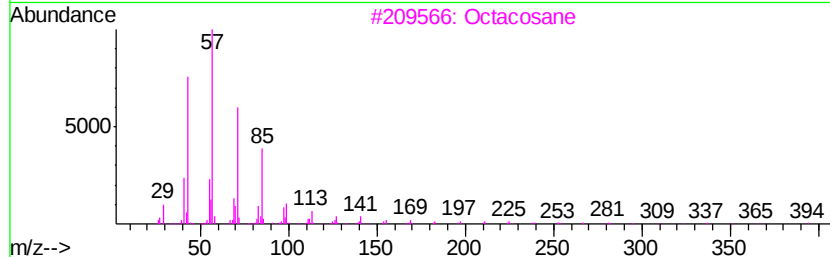
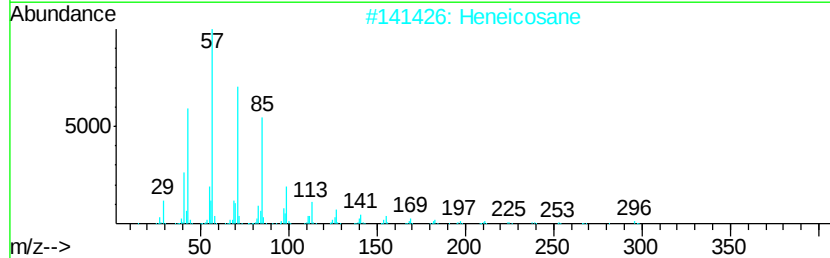
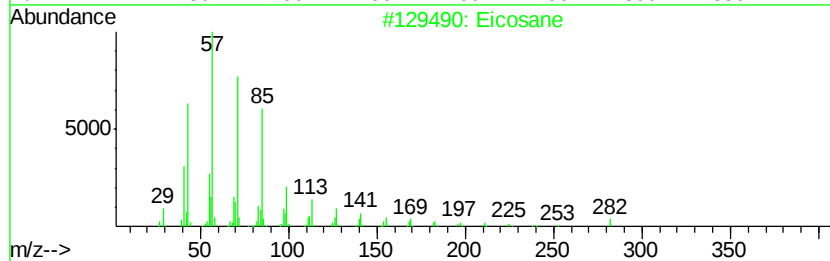
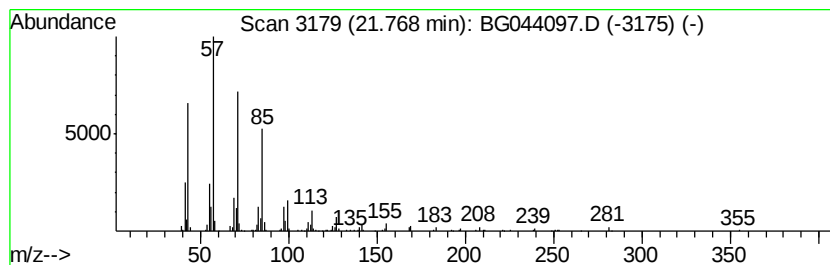
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Peak Number 5 Eicosane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.77	3.14 ng	210495	Chrysene-d12	21.56

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane		282	C20H42	000112-95-8	97
2	Heneicosane		296	C21H44	000629-94-7	91
3	Octacosane		394	C28H58	000630-02-4	91
4	Tetracosane		338	C24H50	000646-31-1	91
5	Eicosane, 7-hexyl-		366	C26H54	055333-99-8	91



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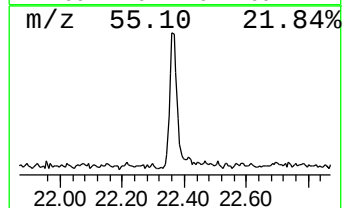
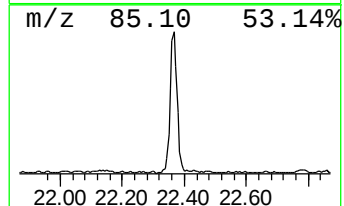
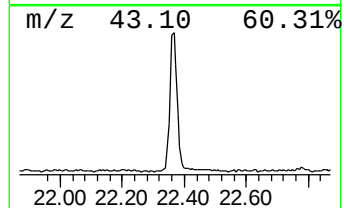
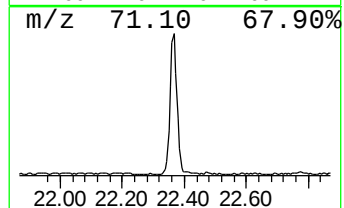
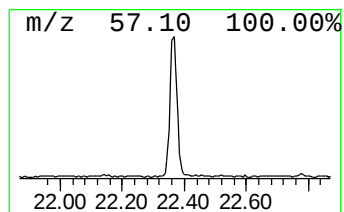
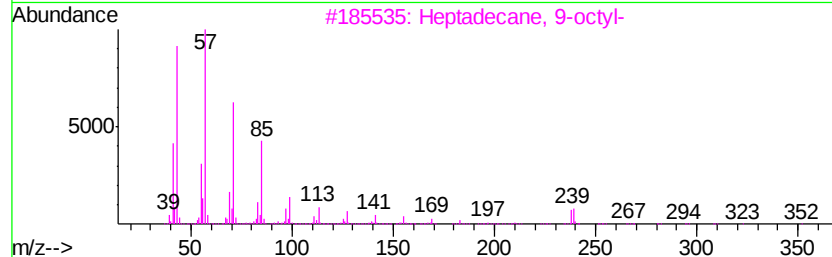
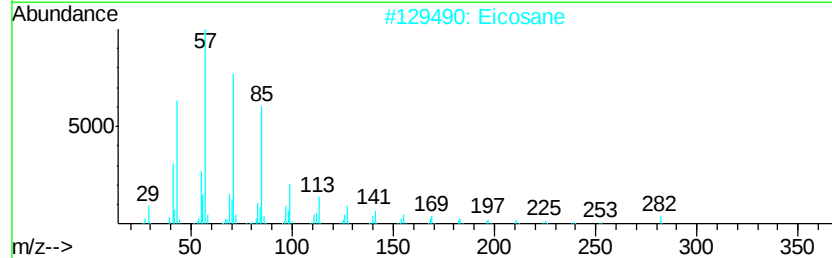
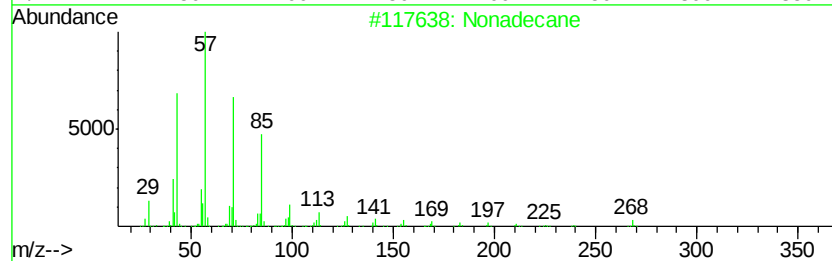
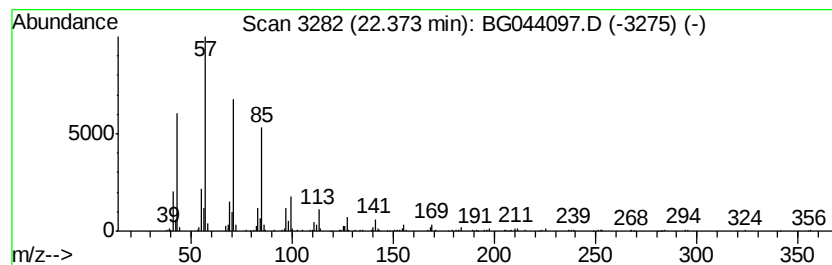
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Peak Number 6 Nonadecane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.37	6.02 ng	404464	Chrysene-d12	21.56

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonadecane		268	C19H40	000629-92-5	94
2	Eicosane		282	C20H42	000112-95-8	94
3	Heptadecane, 9-octyl-		352	C25H52	007225-64-1	91
4	Hexacosane		366	C26H54	000630-01-3	91
5	Tetracosane		338	C24H50	000646-31-1	91



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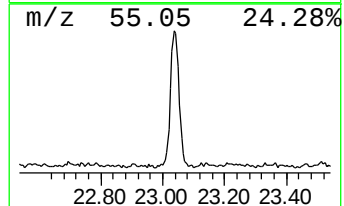
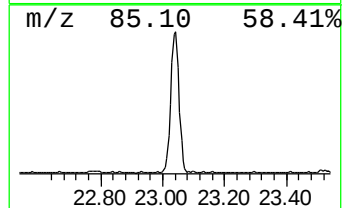
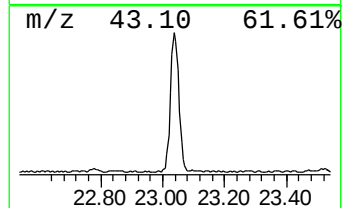
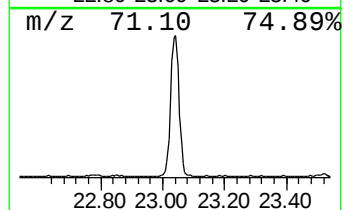
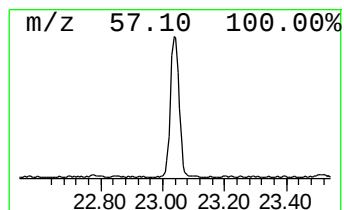
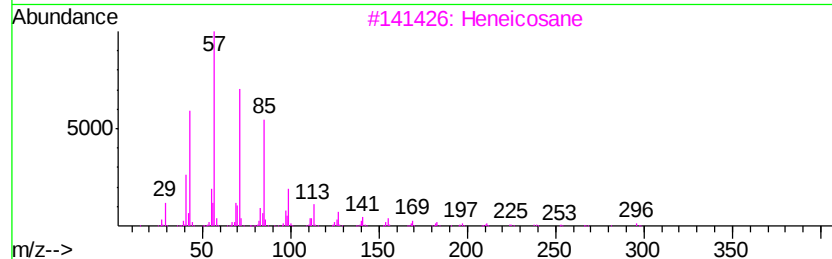
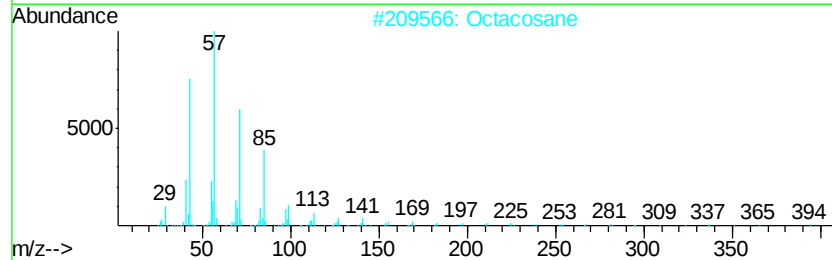
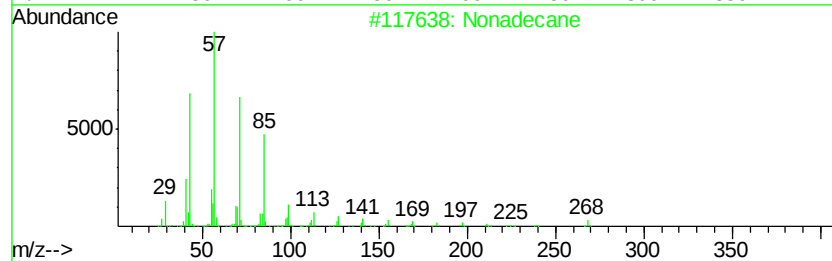
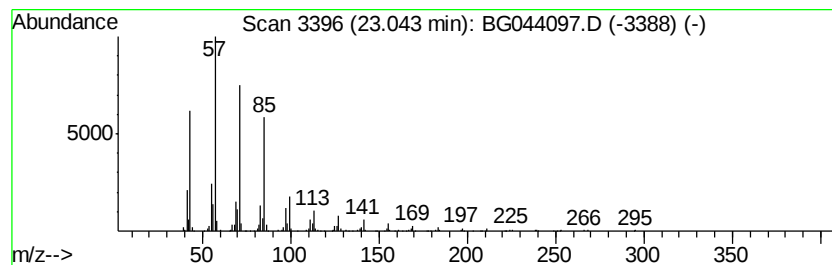
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 Octacosane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.04	8.17 ng	548171	Chrysene-d12	21.56

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonadecane		268	C19H40	000629-92-5	93
2	Octacosane		394	C28H58	000630-02-4	91
3	Heneicosane		296	C21H44	000629-94-7	91
4	Hentriacontane		437	C31H64	000630-04-6	91
5	Heptacosane		380	C27H56	000593-49-7	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG012020\
Data File : BG044097.D
Acq On : 20 Jan 2020 14:27
Operator : JU
Sample : L1176-03
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EP-3

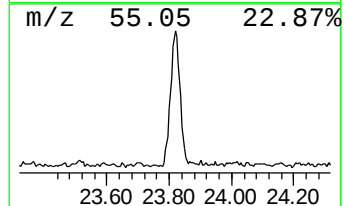
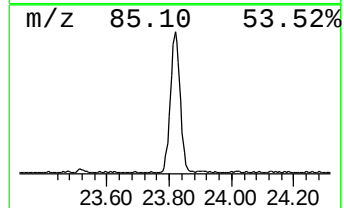
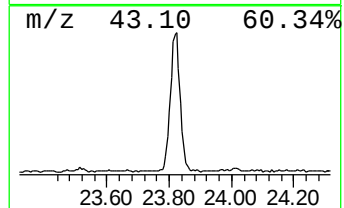
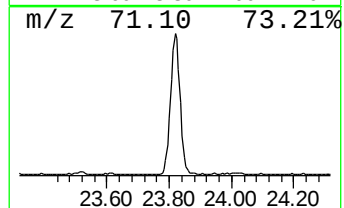
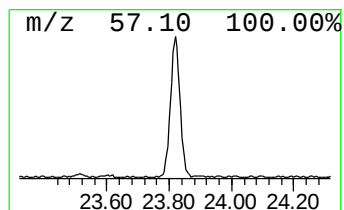
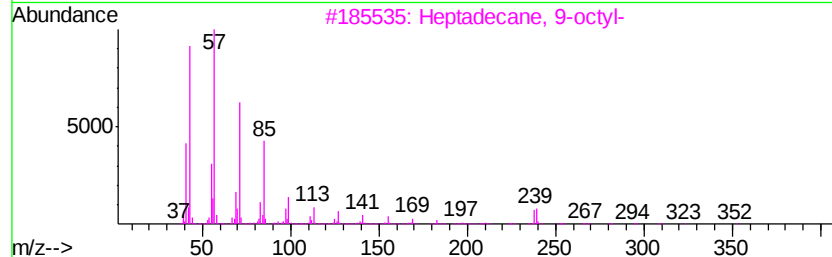
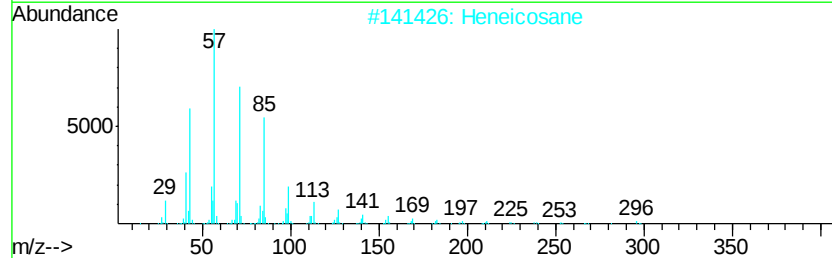
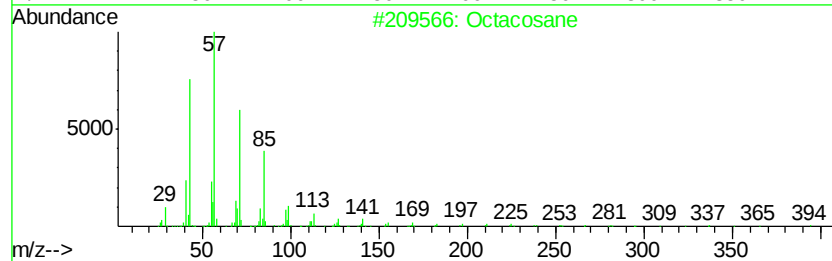
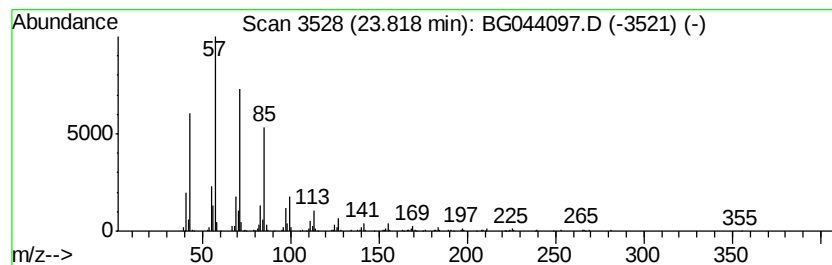
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Heptadecane, 9-octyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.82	8.59 ng	586603	Perylene-d12	24.67

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octacosane	394	C28H58	000630-02-4	91
2		Heneicosane	296	C21H44	000629-94-7	91
3		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
4		Tetracosane	338	C24H50	000646-31-1	91
5		Heneicosane, 11-(1-ethylpropyl)-	366	C26H54	055282-11-6	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG012020\
Data File : BG044097.D
Acq On : 20 Jan 2020 14:27
Operator : JU
Sample : L1176-03
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EP-3

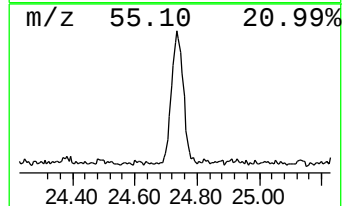
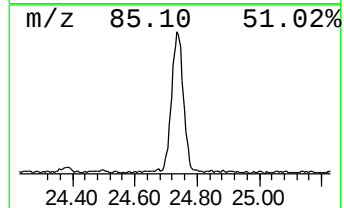
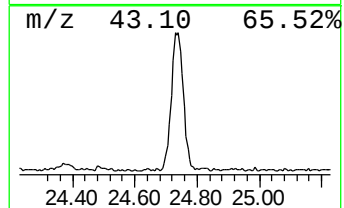
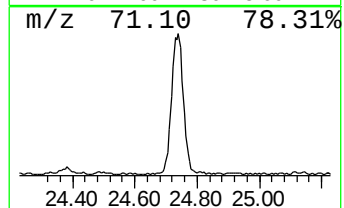
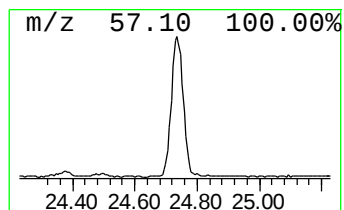
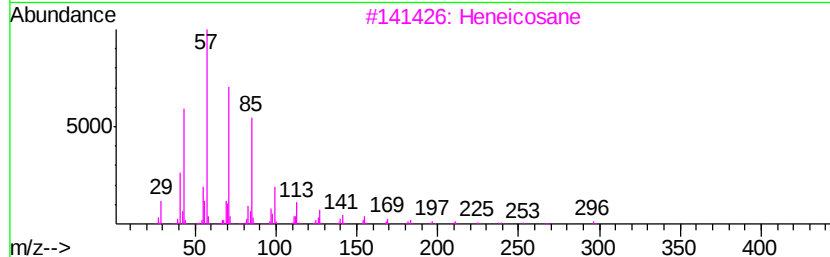
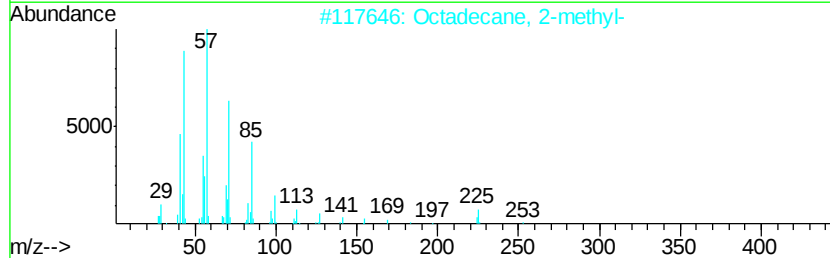
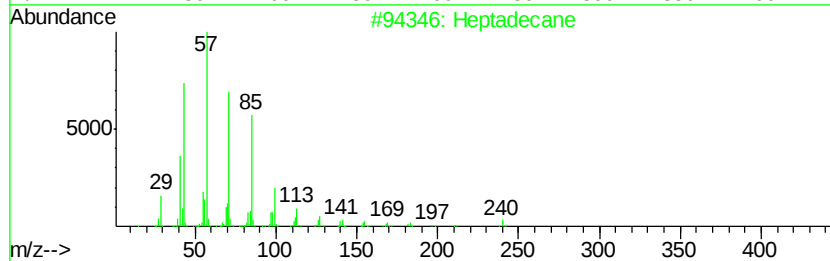
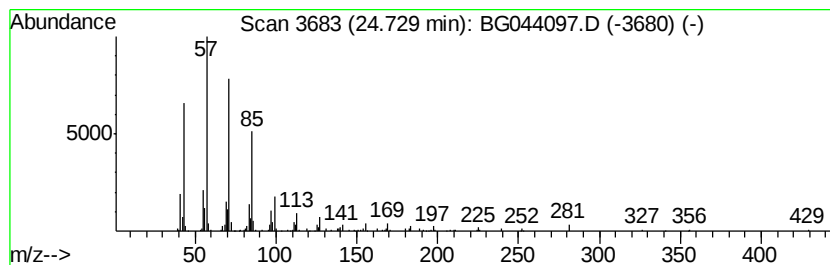
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Heptadecane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.73	7.80 ng	532416	Perylene-d12	24.67

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane		240	C17H36	000629-78-7	87
2	Octadecane, 2-methyl-		268	C19H40	001560-88-9	86
3	Heneicosane		296	C21H44	000629-94-7	86
4	Tridecane, 6-propyl-		226	C16H34	055045-10-8	83
5	Tetracosane		338	C24H50	000646-31-1	83



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG012020\
Data File : BG044097.D
Acq On : 20 Jan 2020 14:27
Operator : JU
Sample : L1176-03
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EP-3

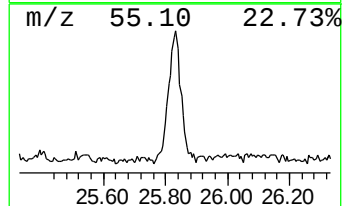
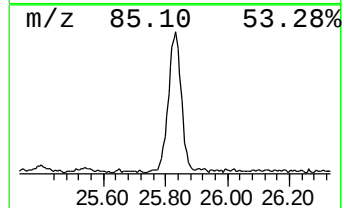
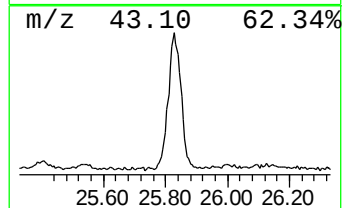
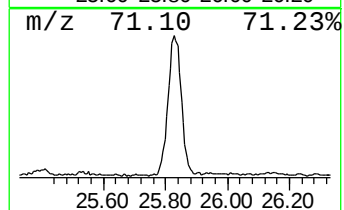
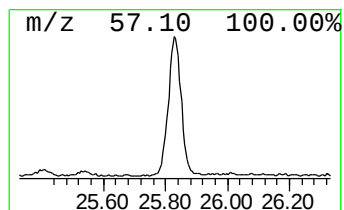
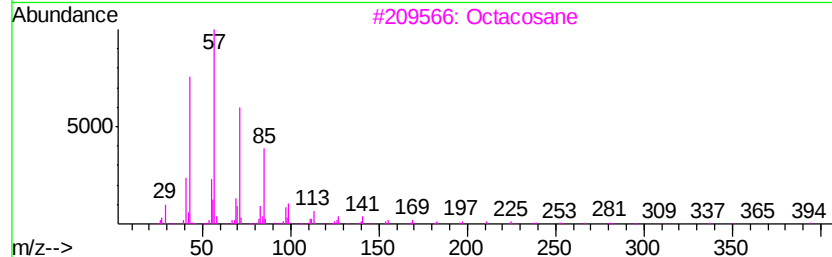
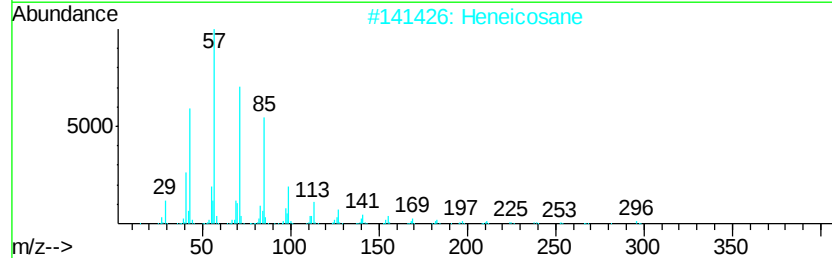
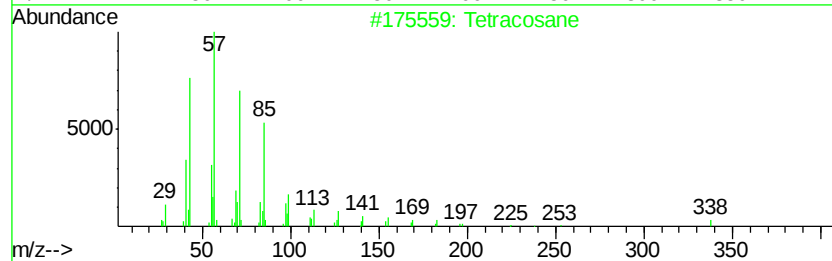
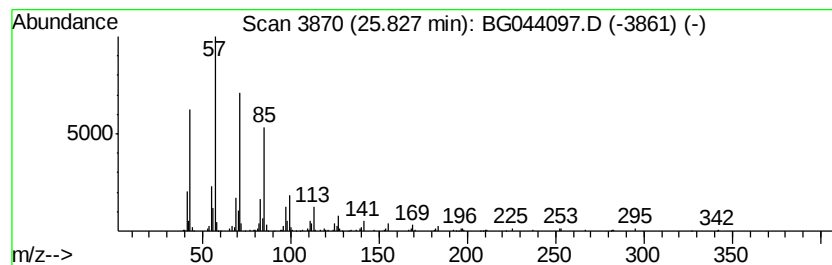
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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.
25.83	6.38 ng	435494	Perylene-d12	24.67

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetracosane	338	C24H50	000646-31-1	91
2			Heneicosane	296	C21H44	000629-94-7	91
3			Octacosane	394	C28H58	000630-02-4	91
4			Heptacosane	380	C27H56	000593-49-7	91
5			Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG012020\
Data File : BG044097.D
Acq On : 20 Jan 2020 14:27
Operator : JU
Sample : L1176-03
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EP-3

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG123019.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--			
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unknown7.47	7.47	73.2	ng	1825900	1	7.93	499130	20.0
Oxalic acid, isob...	21.24	5.2	ng	348736	5	21.56	1342670	20.0
Eicosane	21.77	3.1	ng	210495	5	21.56	1342670	20.0
Nonadecane	22.37	6.0	ng	404464	5	21.56	1342670	20.0
Octacosane	23.04	8.2	ng	548171	5	21.56	1342670	20.0
Heptadecane, 9-oc...	23.82	8.6	ng	586603	6	24.67	1365990	20.0
Heptadecane	24.73	7.8	ng	532416	6	24.67	1365990	20.0
Tetracosane	25.83	6.4	ng	435494	6	24.67	1365990	20.0