

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG022819\
 Data File : BG039673.D
 Acq On : 28 Feb 2019 23:32
 Operator : JU/SJ
 Sample : K1643-03
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 FLOOR-SLAB

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-BG012919.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	3.220	16	22	31	rBV2	83020	169845	3.96%	0.612%
2	3.296	31	35	43	rVB	54958	79932	1.87%	0.288%
3	4.635	258	263	273	rBV	138596	211930	4.95%	0.764%
4	5.246	360	367	376	rBV	267592	406602	9.49%	1.466%
5	5.710	439	446	460	rVB	1031535	1621480	37.85%	5.846%
6	7.314	709	719	731	rBV	1224844	2113927	49.35%	7.622%
7	7.696	775	784	798	rBV	1047049	1856280	43.33%	6.693%
8	8.166	856	864	874	rBV	208083	358515	8.37%	1.293%
9	8.489	907	919	927	rBV	856352	1482369	34.60%	5.345%
10	9.341	1054	1064	1074	rBV	724849	1321740	30.85%	4.766%
11	10.986	1336	1344	1353	rVB	273576	502485	11.73%	1.812%
12	13.412	1749	1757	1764	rBV	2031410	3242262	75.68%	11.690%
13	14.228	1891	1896	1901	rBV	99924	140712	3.28%	0.507%
14	14.786	1985	1991	2002	rBV2	458071	758897	17.72%	2.736%
15	16.273	2237	2244	2259	rBV	1340355	2083634	48.64%	7.513%
16	17.530	2452	2458	2470	rBV2	601623	975808	22.78%	3.518%
17	18.381	2598	2603	2609	rBV2	44354	74358	1.74%	0.268%
18	18.464	2613	2617	2625	rVB	49429	67726	1.58%	0.244%
19	19.227	2742	2747	2754	rBV	109676	187236	4.37%	0.675%
20	19.292	2754	2758	2765	rVB	106363	141201	3.30%	0.509%
21	19.768	2835	2839	2843	rVB	42062	52682	1.23%	0.190%
22	19.862	2851	2855	2859	rVB	124064	142041	3.32%	0.512%
23	20.126	2894	2900	2906	rBV	3203559	4283907	100.00%	15.446%
24	20.390	2942	2945	2949	rBV	114503	129840	3.03%	0.468%
25	20.890	3026	3030	3034	rVB	135986	149634	3.49%	0.540%
26	21.407	3114	3118	3123	rVB	188114	236567	5.52%	0.853%
27	21.665	3157	3162	3171	rVB	324304	459063	10.72%	1.655%
28	21.818	3182	3188	3199	rVB2	555765	977047	22.81%	3.523%
29	21.971	3209	3214	3219	rVB	149072	215733	5.04%	0.778%
30	22.599	3315	3321	3328	rBV	192579	315064	7.35%	1.136%
31	22.993	3384	3388	3395	rVB3	62048	112780	2.63%	0.407%
32	23.316	3437	3443	3451	rVB	183087	354460	8.27%	1.278%
33	24.156	3578	3586	3594	rBV2	226040	503216	11.75%	1.814%
34	25.149	3747	3755	3756	rBV2	161298	302432	7.06%	1.090%

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

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
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35	25.196	3758	3763	3777	rVB2	305802	906099	21.15%	3.267%
36	26.330	3946	3956	3969	rVB3	141478	458584	10.70%	1.653%
37	27.757	4186	4199	4210	rBV4	55660	223810	5.22%	0.807%
38	29.478	4486	4492	4504	rVB7	30338	114759	2.68%	0.414%

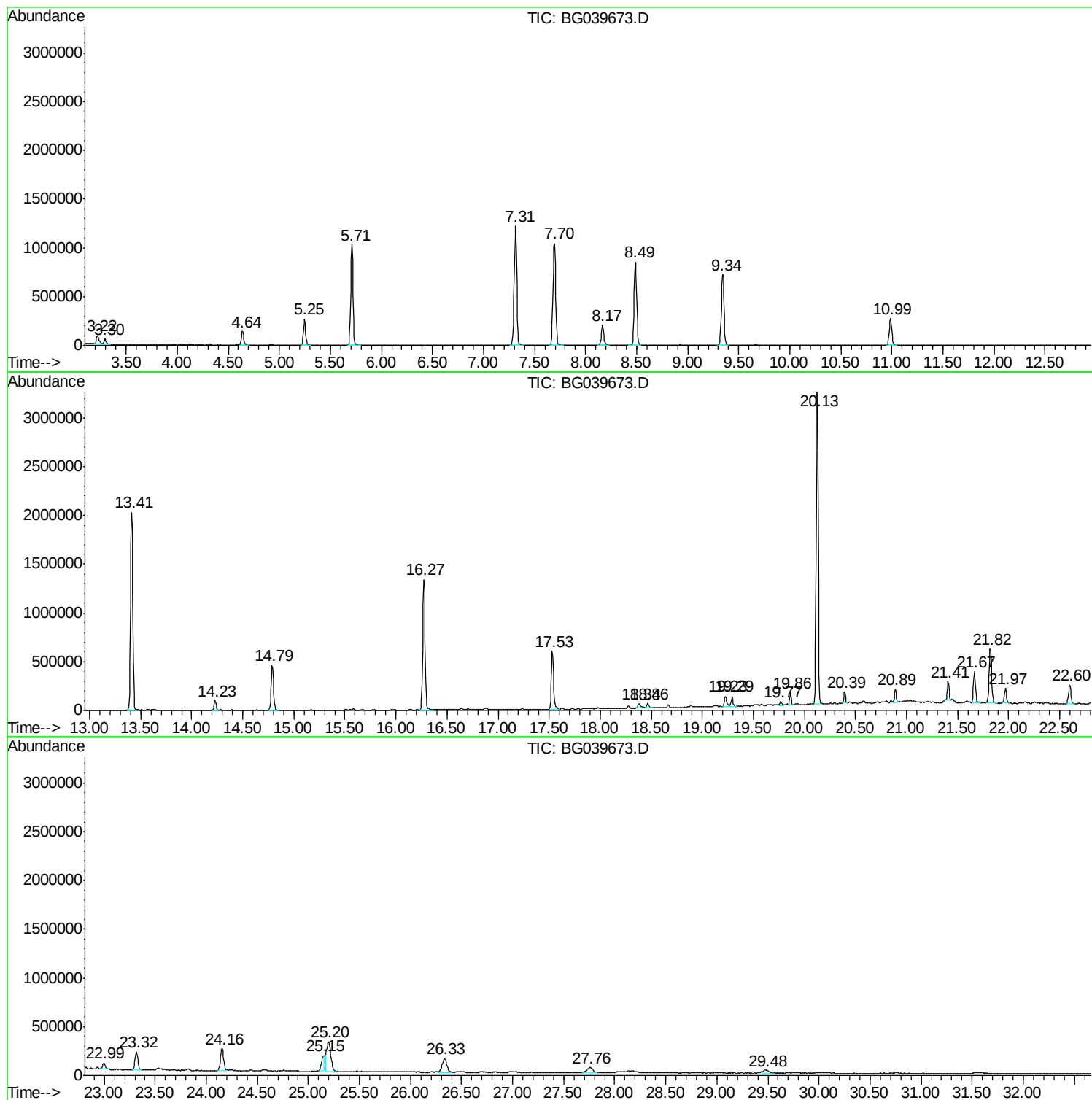
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG012919.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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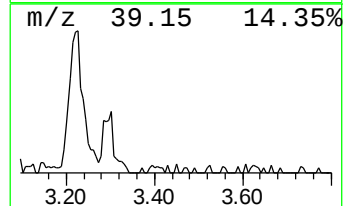
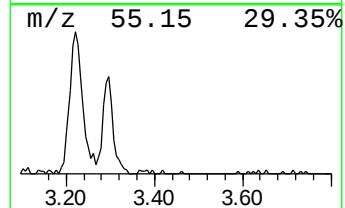
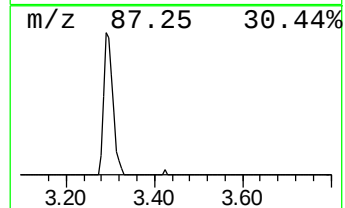
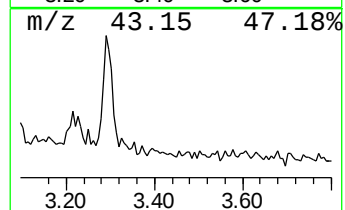
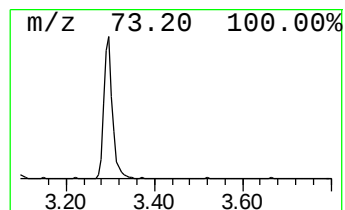
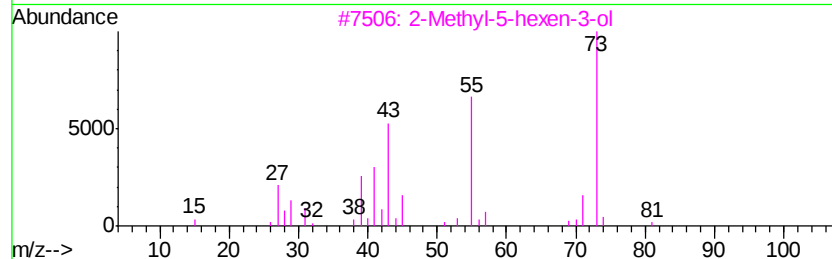
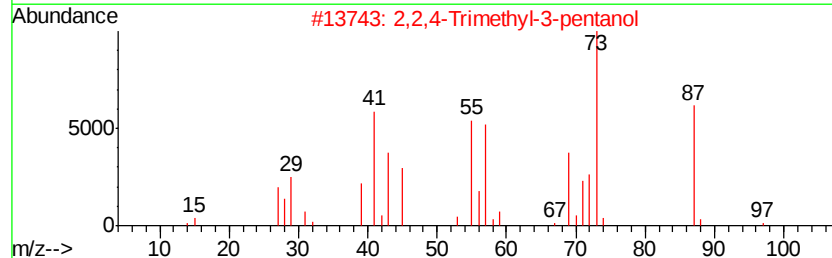
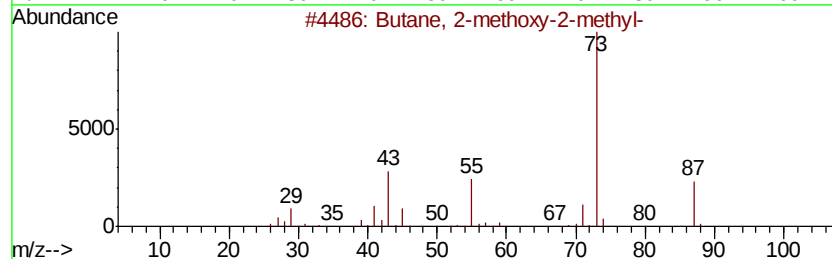
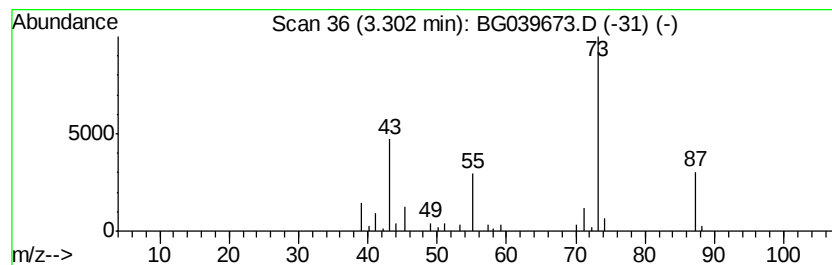
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 Butane, 2-methoxy-2-methyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.30	4.46 ng	79932	1,4-Dichlorobenzene-d4	8.17

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2			2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	56
3			2-Methyl-5-hexen-3-ol	114	C7H14O	032815-70-6	9
4			3-Pentanol, 2,4-dimethyl-	116	C7H16O	000600-36-2	9
5			Silane, ethyltrimethyl-	102	C5H14Si	003439-38-1	9



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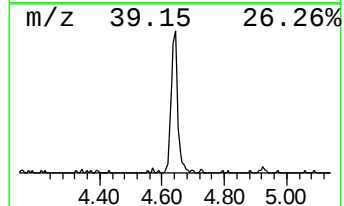
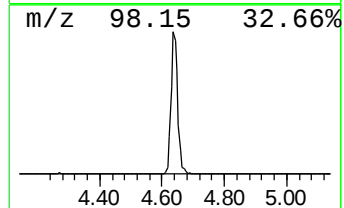
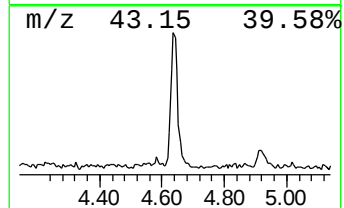
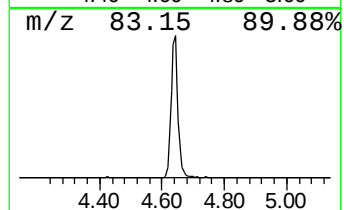
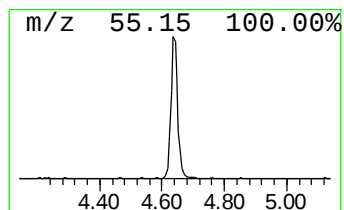
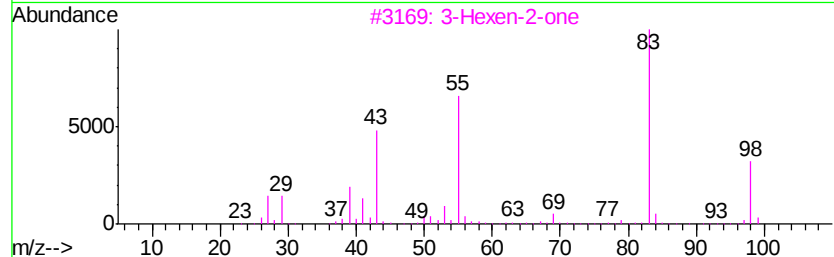
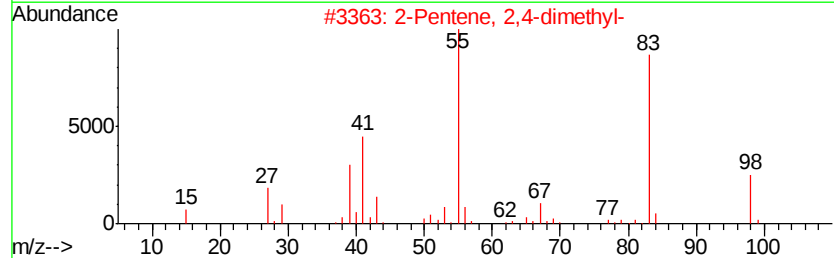
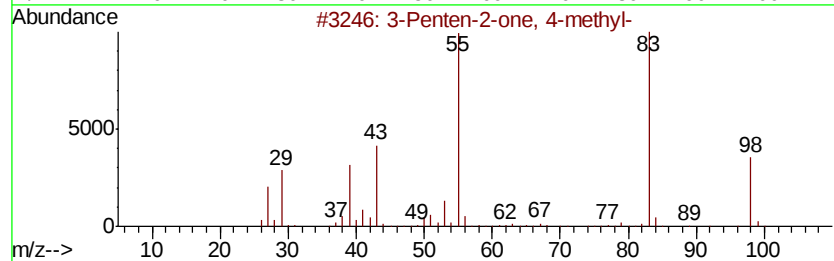
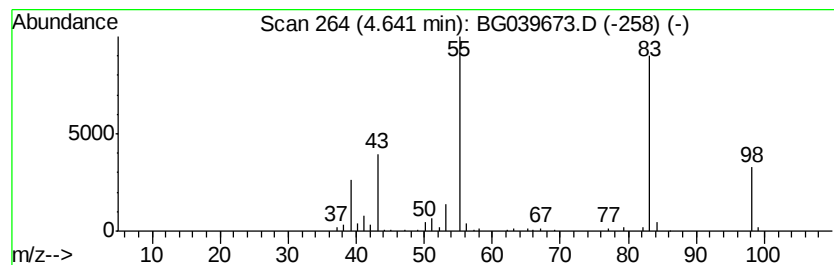
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 3-Penten-2-one, 4-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.64	11.82 ng	211930	1,4-Dichlorobenzene-d4	8.17

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Penten-2-one, 4-methyl-	98	C6H10O	000141-79-7	90
2			2-Pentene, 2,4-dimethyl-	98	C7H14	000625-65-0	86
3			3-Hexen-2-one	98	C6H10O	000763-93-9	86
4			2-Pentene, 3,4-dimethyl-	98	C7H14	024910-63-2	78
5			2-Pentene, 3,4-dimethyl-, (Z)-	98	C7H14	004914-91-4	78



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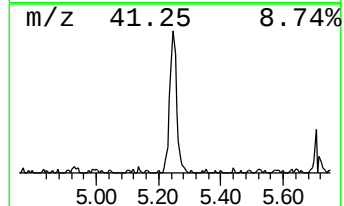
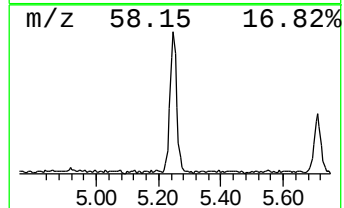
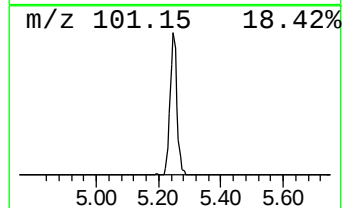
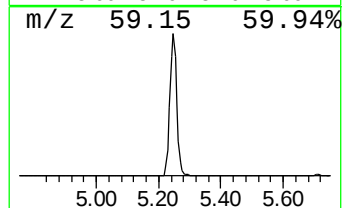
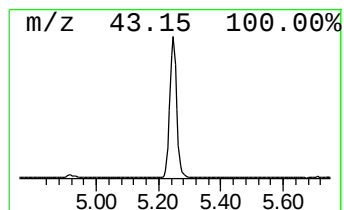
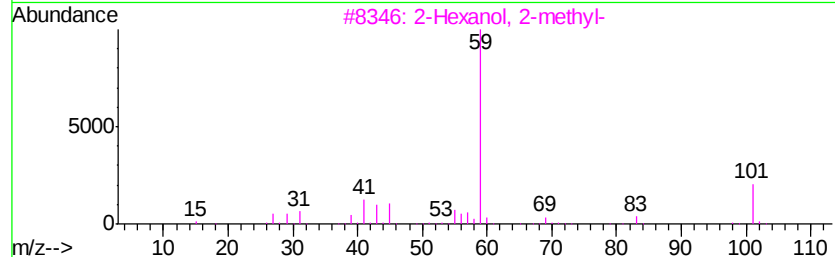
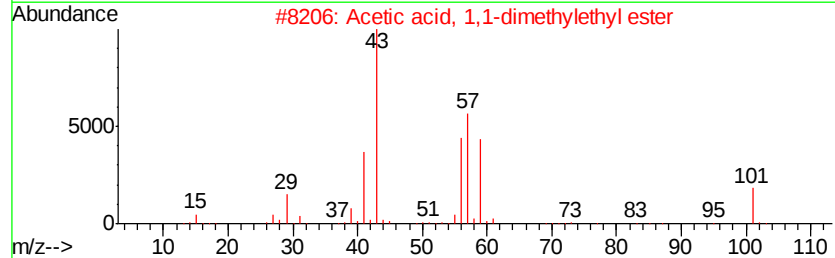
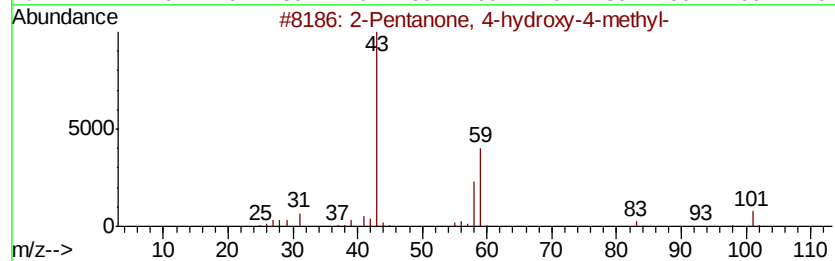
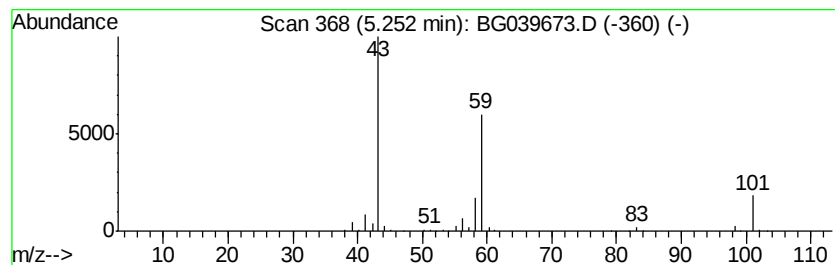
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.25	22.68 ng	406602	1,4-Dichlorobenzene-d4	8.17

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2		Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	38
3		2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	33
4		Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	23
5		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	10



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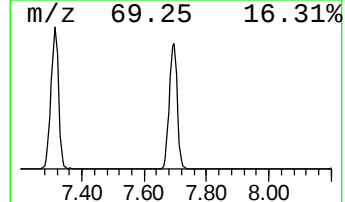
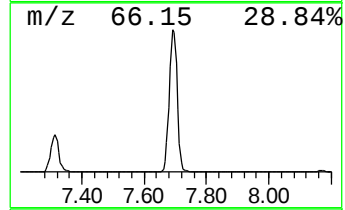
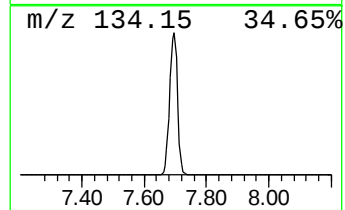
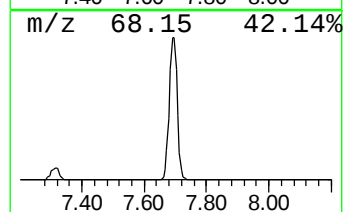
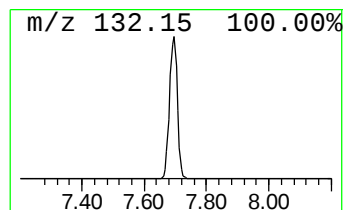
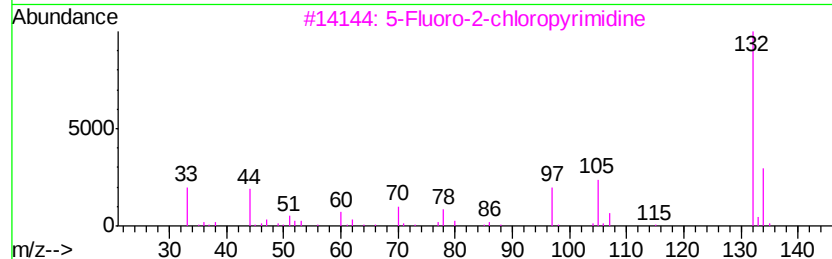
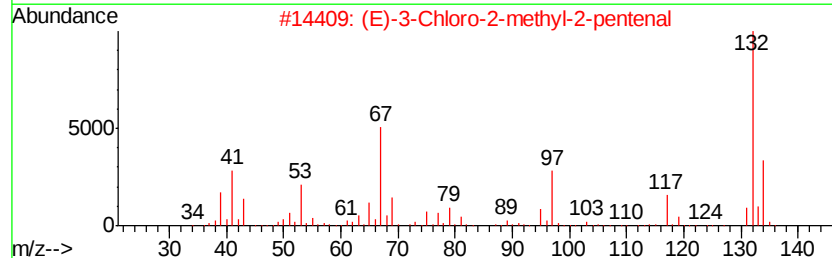
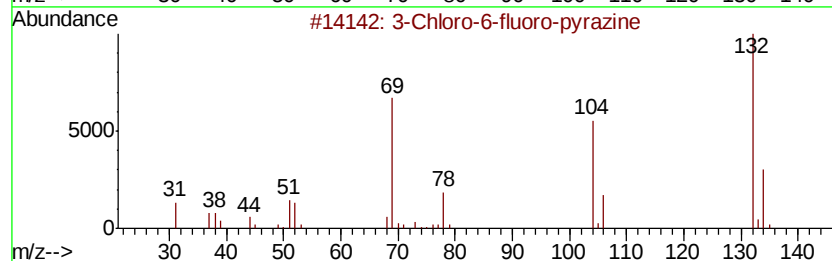
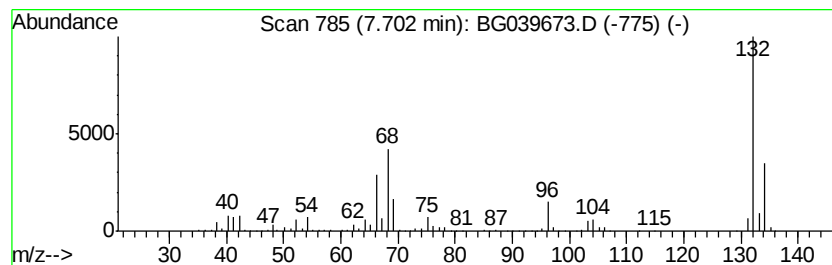
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Peak Number 5 unknown7.70 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.70	103.55 ng	1856280	1,4-Dichlorobenzene-d4	8.17

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		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
4		(5-Methyl-2-pyridyl)acetonitrile	132	C8H8N2	1000241-93-9	10
5		Xenon	132	Xe	007440-63-3	9



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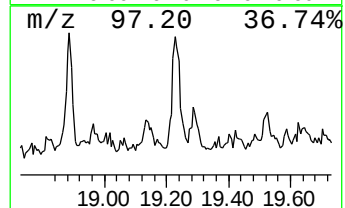
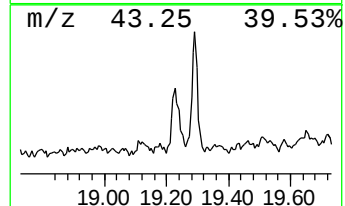
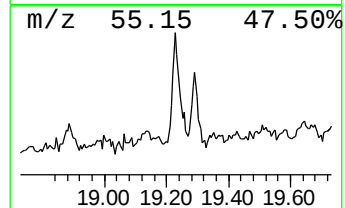
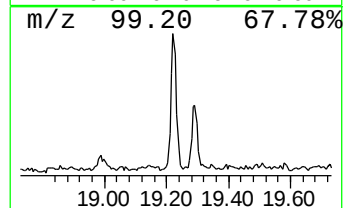
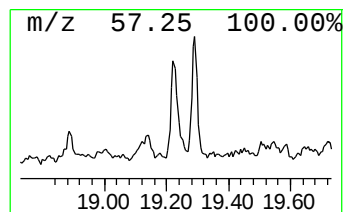
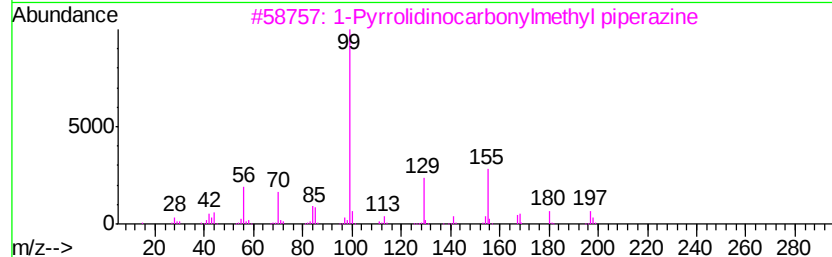
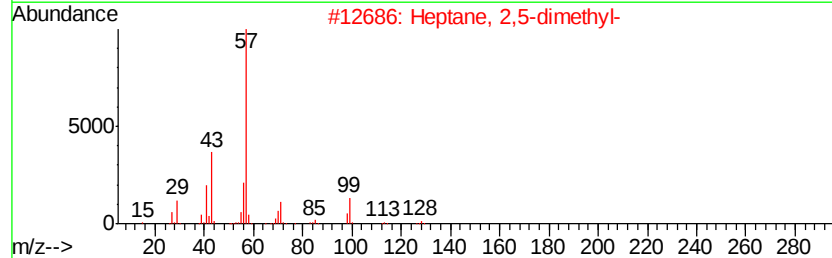
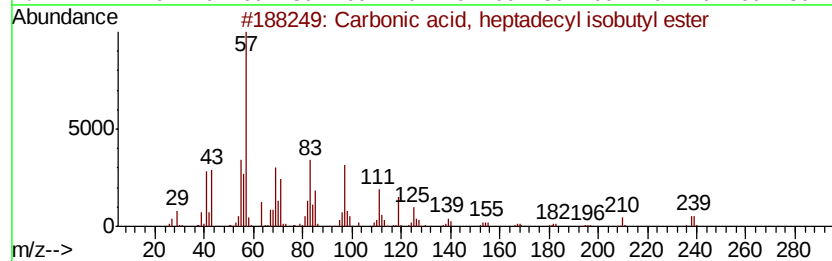
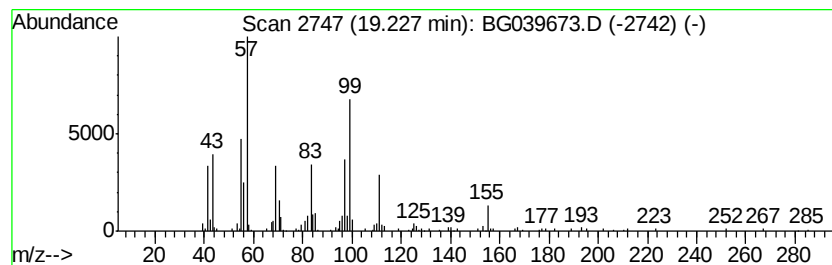
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Peak Number 7 unknown19.23 Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.23	3.84 ng	187236	Phenanthrene-d10	17.53

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Carbonic acid, heptadecyl isobut...	356	C22H44O3	1000314-61-4	43
2		Heptane, 2,5-dimethyl-	128	C9H20	002216-30-0	43
3		1-Pyrrolidinocarbonylmethyl pipe...	197	C10H19N3O	039890-45-4	38
4		Methylphosphonic acid, fluoroanh...	238	C11H24F02P	211192-71-1	35
5		2-Ethyl-1-hexanol, heptafluorobu...	326	C12H17F7O2	1000365-16-0	27



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Acq On : 28 Feb 2019 23:32
Operator : JU/SJ
Sample : K1643-03
Misc :
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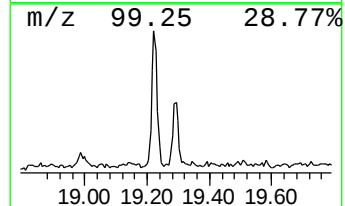
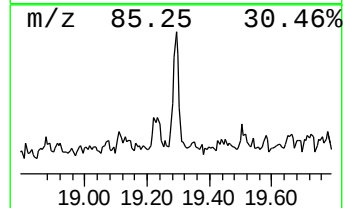
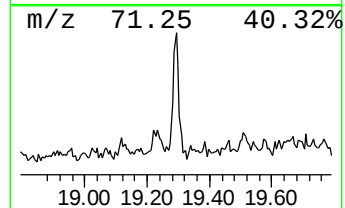
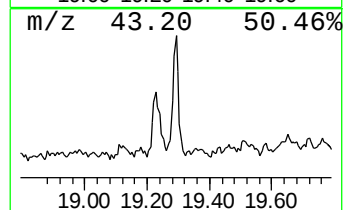
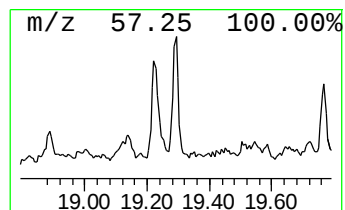
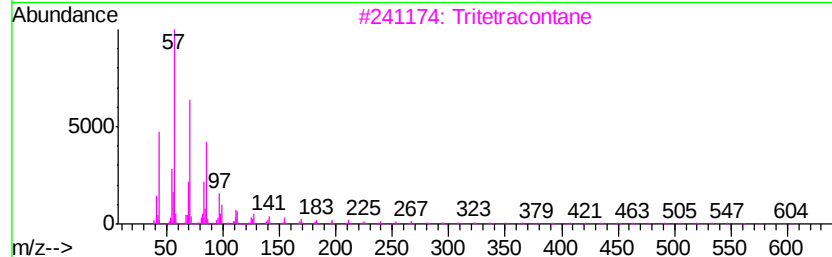
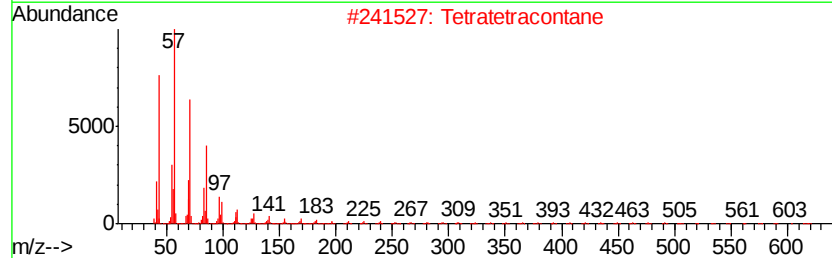
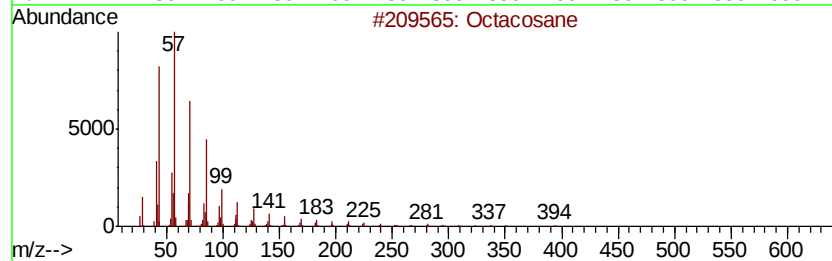
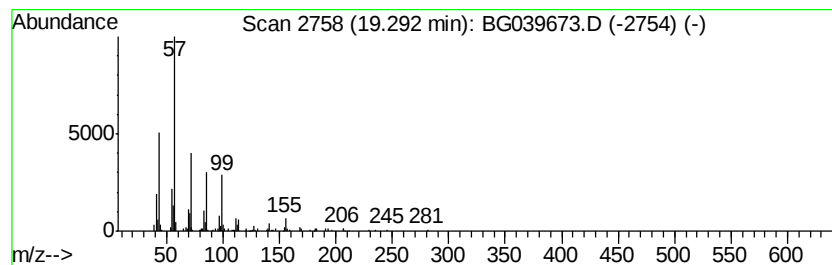
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Octacosane Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.29	2.89 ng	141201	Phenanthrene-d10	17.53

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octacosane		394	C28H58	000630-02-4	64
2	Tetratetracontane		619	C44H90	007098-22-8	58
3	Tritetracontane		605	C43H88	007098-21-7	58
4	Sulfurous acid, butyl decyl ester		278	C14H30O3S	1000309-17-7	58
5	Pentacosane		352	C25H52	000629-99-2	58



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG022819\
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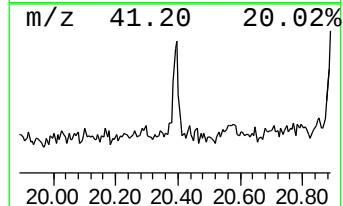
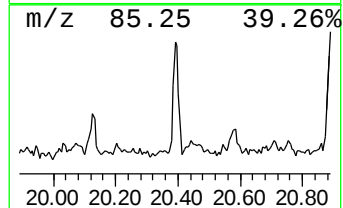
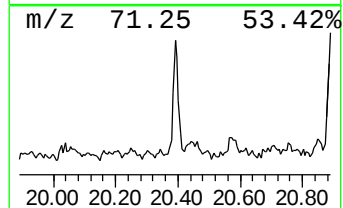
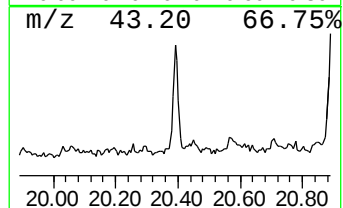
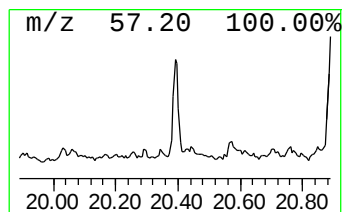
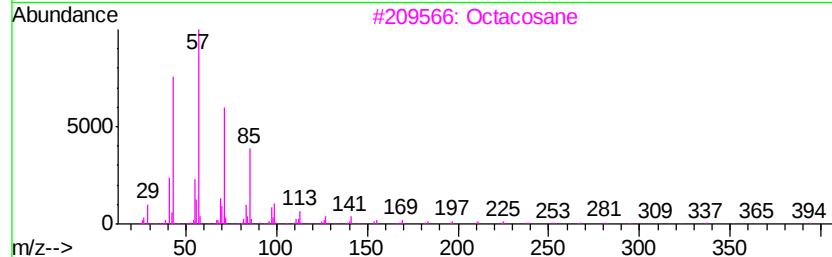
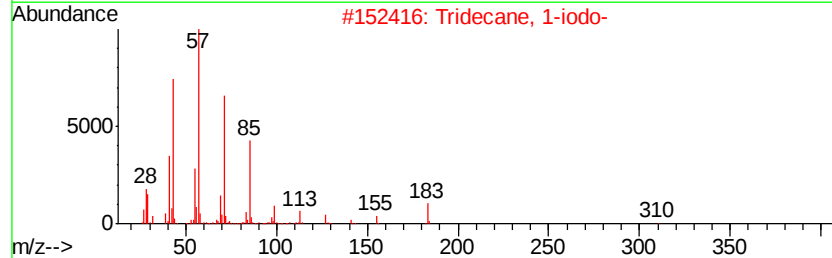
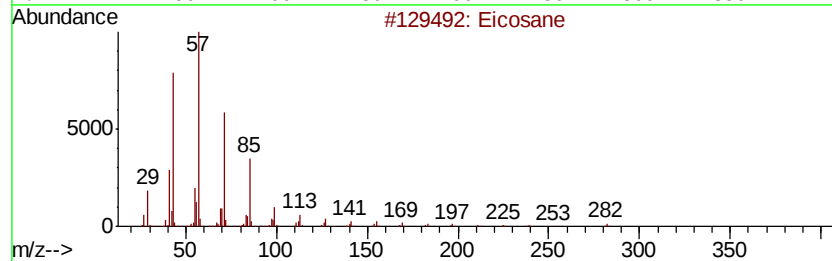
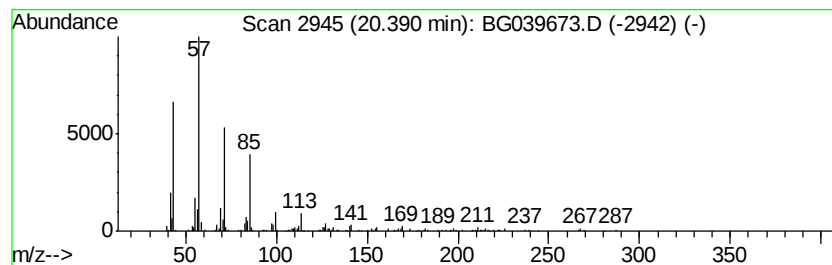
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TIC Library : C:\Database\NIST11.L
 TIC Integration Parameters: LSCINT.P

 Peak Number 10 Eicosane Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.39	2.66 ng	129840	Chrysene-d12	21.82

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane	282	C20H42	000112-95-8	90
2		Tridecane, 1-iodo-	310	C13H27I	035599-77-0	90
3		Octacosane	394	C28H58	000630-02-4	90
4		Hexadecane	226	C16H34	000544-76-3	87
5		Tridecane, 7-hexyl-	268	C19H40	007225-66-3	86



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG022819\
 Data File : BG039673.D
 Acq On : 28 Feb 2019 23:32
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Instrument :
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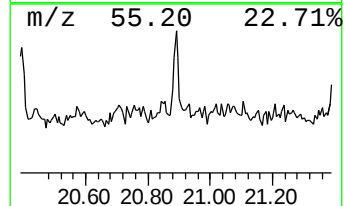
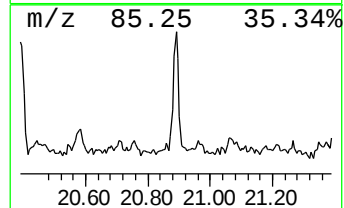
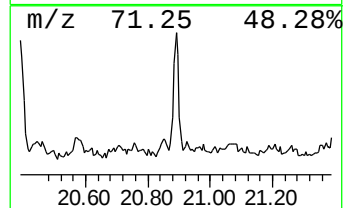
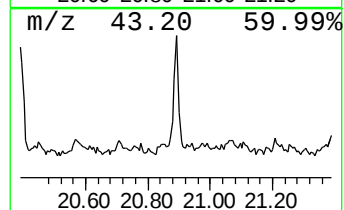
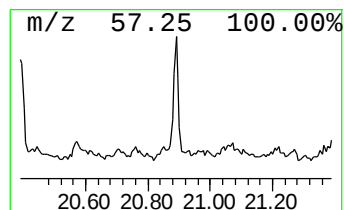
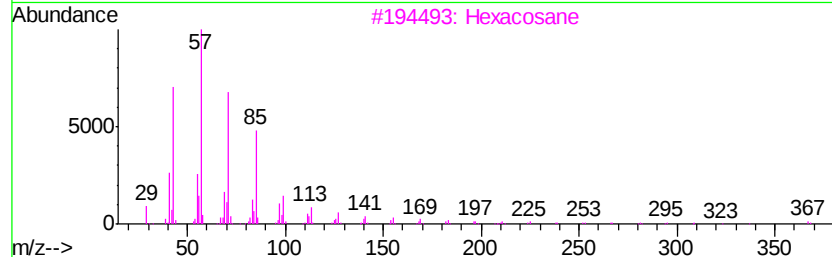
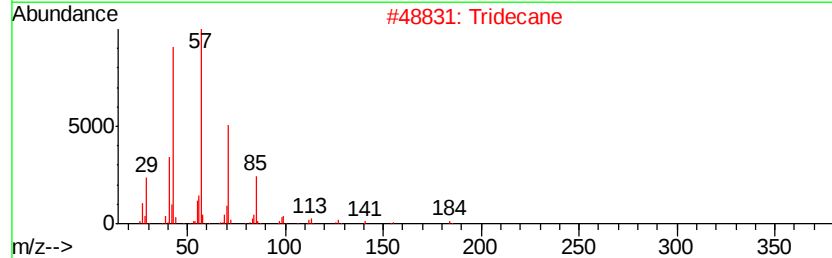
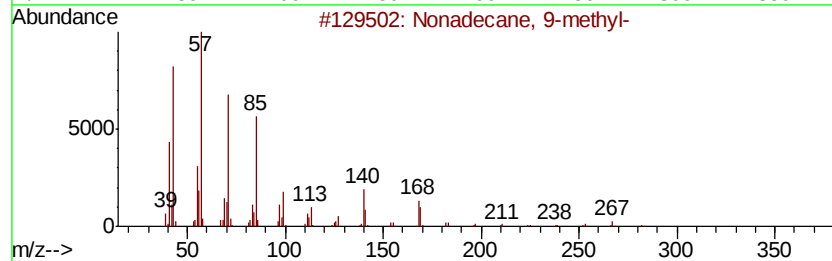
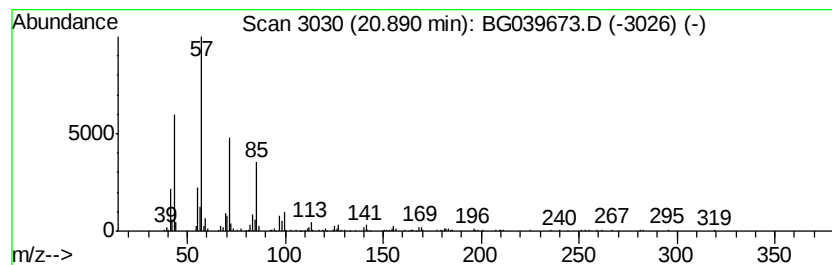
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TIC Library : C:\Database\NIST11.L
 TIC Integration Parameters: LSCINT.P

 Peak Number 11 Nonadecane, 9-methyl- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.89	3.06 ng	149634	Chrysene-d12	21.82

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Nonadecane, 9-methyl-	282	C20H42	013287-24-6	93
2		Tridecane	184	C13H28	000629-50-5	90
3		Hexacosane	366	C26H54	000630-01-3	90
4		Eicosane	282	C20H42	000112-95-8	90
5		Hexadecane	226	C16H34	000544-76-3	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG022819\
Data File : BG039673.D
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Sample : K1643-03
Misc :
ALS Vial : 15 Sample Multiplier: 1

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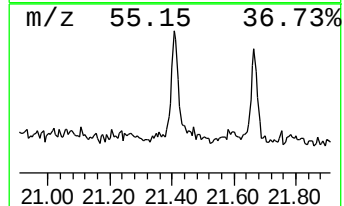
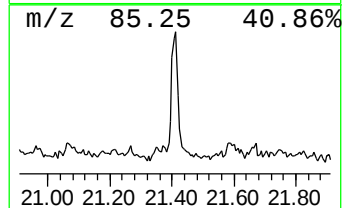
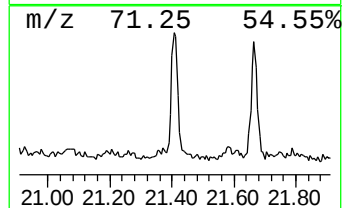
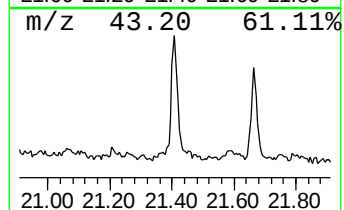
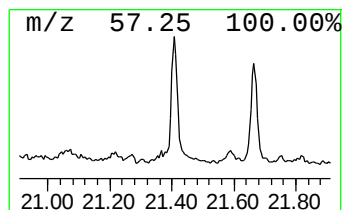
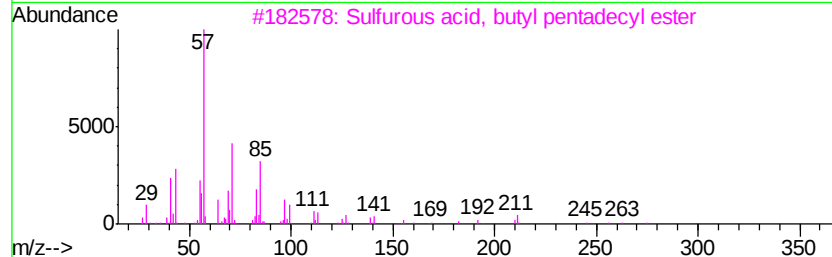
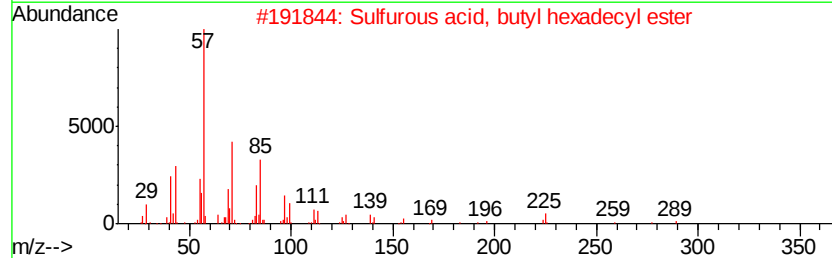
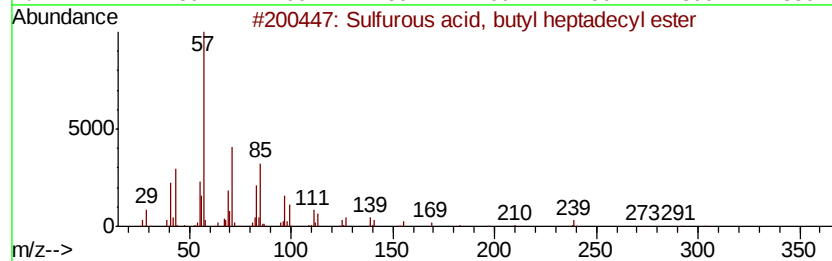
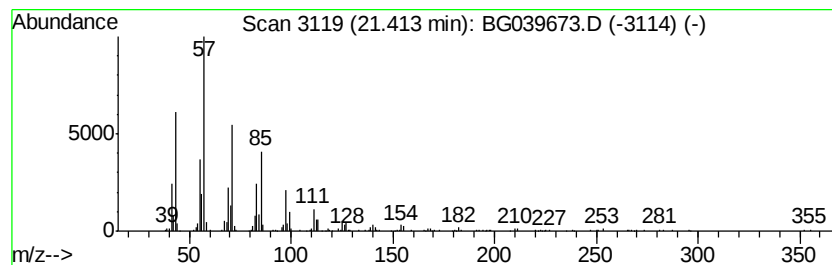
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Sulfurous acid, butyl hepta... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.41	4.84 ng	236567	Chrysene-d12	21.82

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Sulfurous acid, butyl heptadecyl...	376	C21H44O3S	1000309-18-4	91
2			Sulfurous acid, butyl hexadecyl ...	362	C20H42O3S	1000309-18-3	90
3			Sulfurous acid, butyl pentadecyl...	348	C19H40O3S	1000309-18-2	90
4			Sulfurous acid, butyl tridecyl e...	320	C17H36O3S	1000309-18-0	86
5			Heptacosane, 1-chloro-	414	C27H55Cl	062016-79-9	86



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG022819\
Data File : BG039673.D
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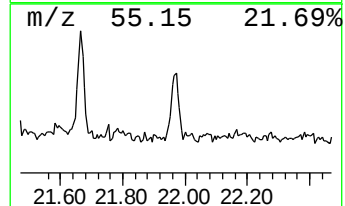
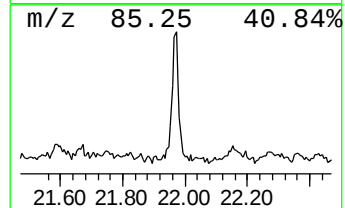
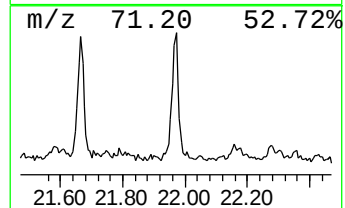
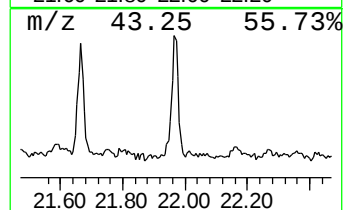
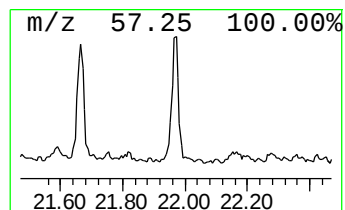
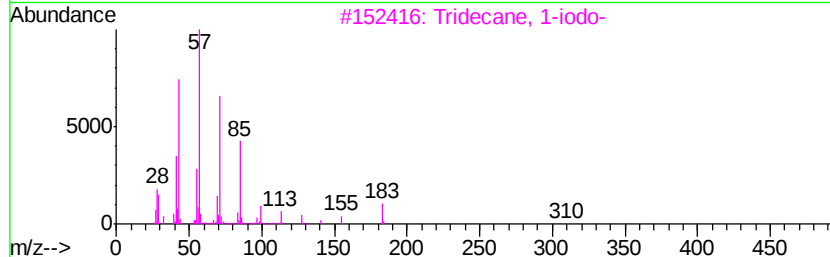
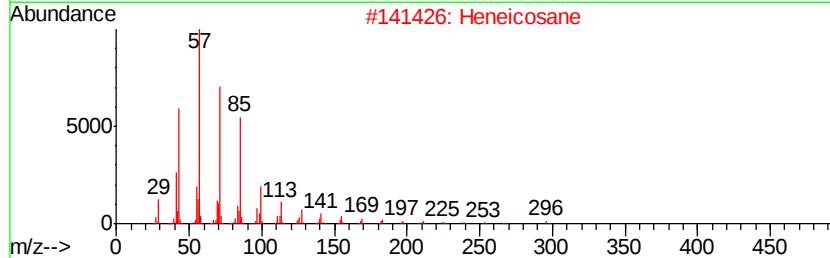
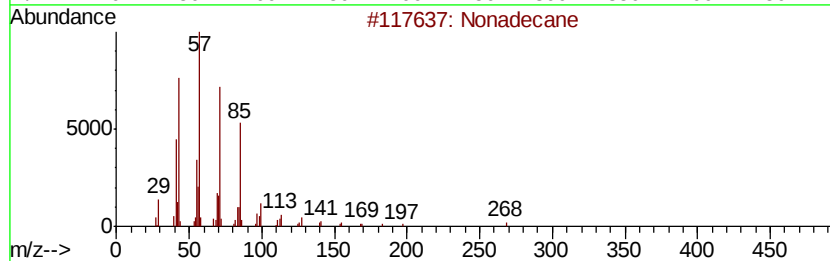
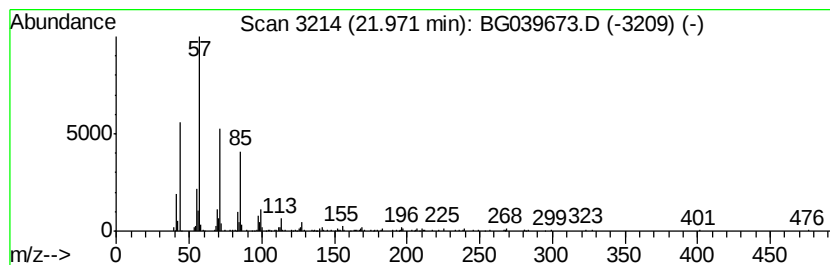
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 13 Nonadecane Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.97	4.42 ng	215733	Chrysene-d12	21.82

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Nonadecane	268	C19H40	000629-92-5	87
2		Heneicosane	296	C21H44	000629-94-7	74
3		Tridecane, 1-iodo-	310	C13H27I	035599-77-0	72
4		Heptadecane	240	C17H36	000629-78-7	72
5		Heptacosane	380	C27H56	000593-49-7	68



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG022819\
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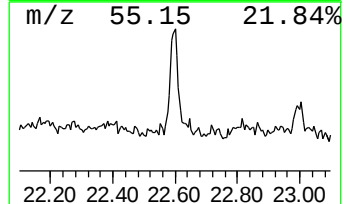
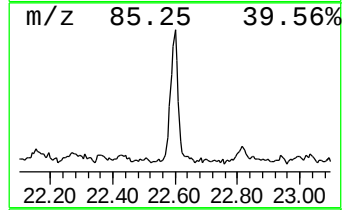
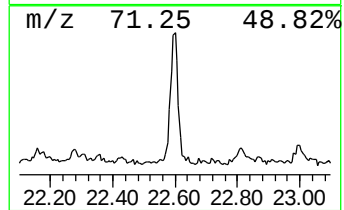
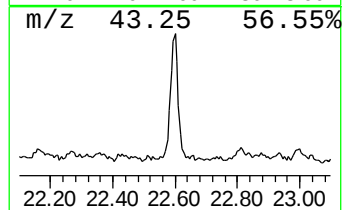
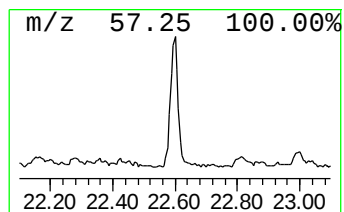
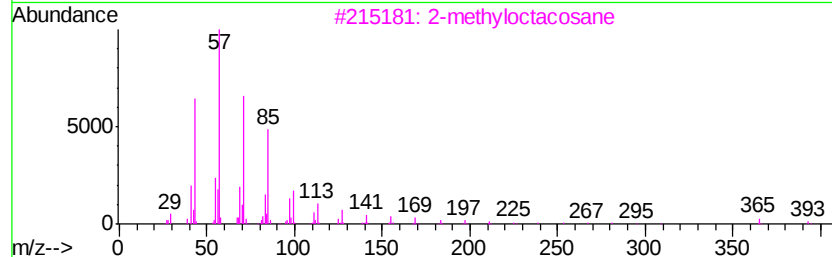
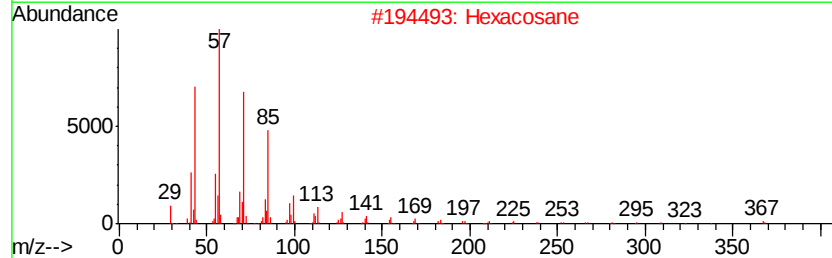
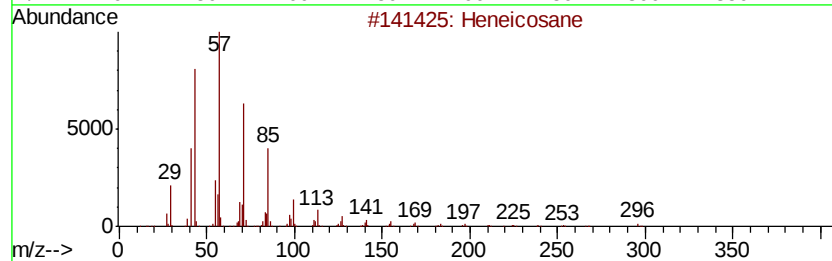
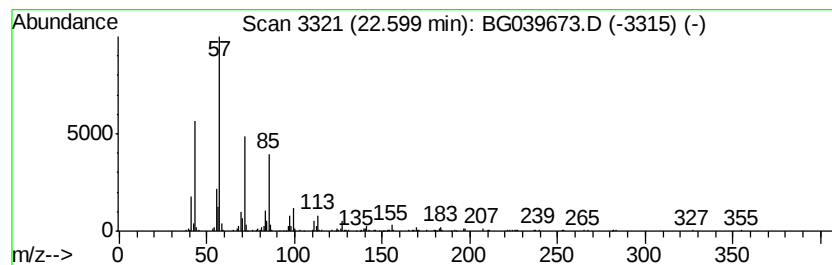
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 14 Heneicosane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.60	6.45 ng	315064	Chrysene-d12	21.82

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosane	296	C21H44	000629-94-7	91
2		Hexacosane	366	C26H54	000630-01-3	90
3		2-methyloctacosane	408	C29H60	1000376-72-8	90
4		Octadecane	254	C18H38	000593-45-3	87
5		Eicosane	282	C20H42	000112-95-8	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG022819\
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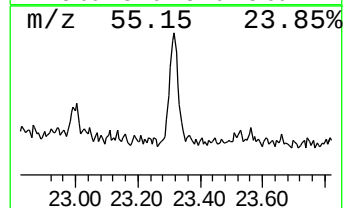
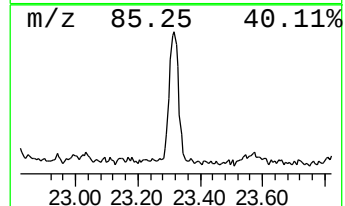
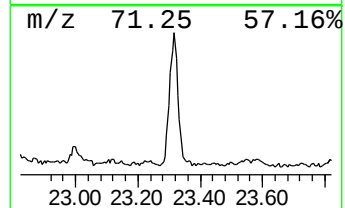
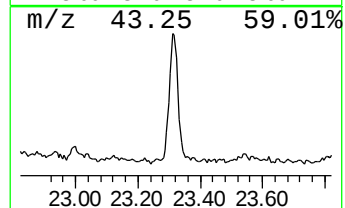
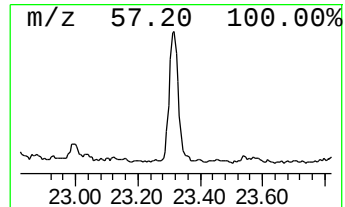
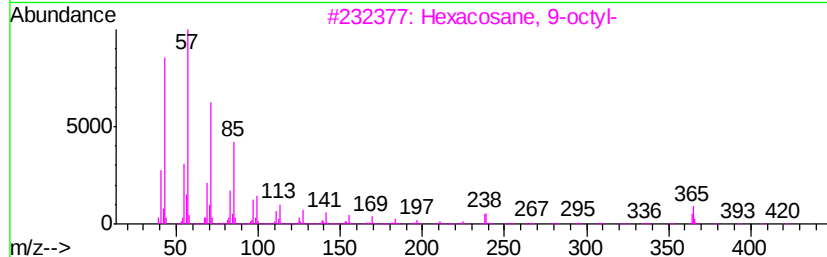
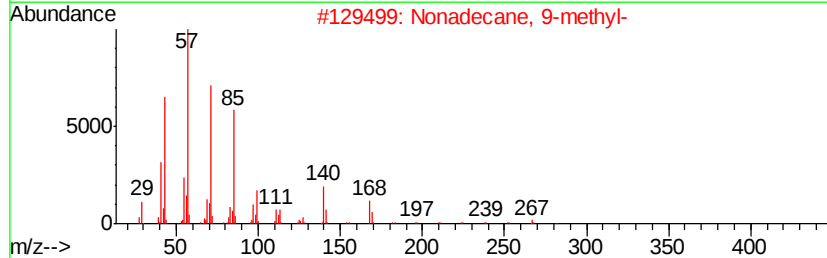
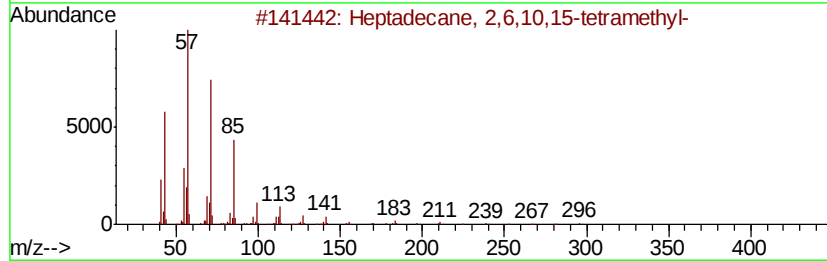
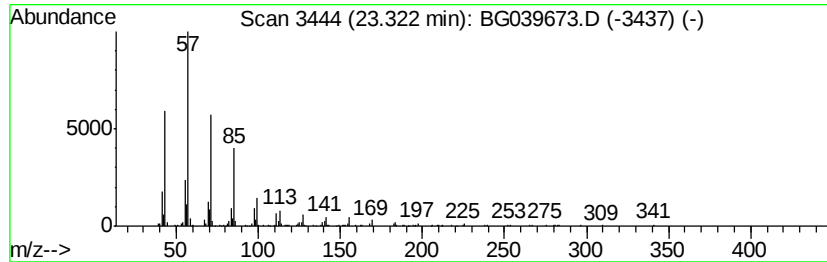
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

 Peak Number 15 Heptadecane, 2,6,10,15-tetr... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.32	7.26 ng	354460	Chrysene-d12	21.82

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	97
2		Nonadecane, 9-methyl-	282	C20H42	013287-24-6	93
3		Hexacosane, 9-octyl-	479	C34H70	055429-83-9	91
4		Hentriacontane	437	C31H64	000630-04-6	91
5		Tetracosane, 11-decyl-	479	C34H70	055429-84-0	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG022819\
Data File : BG039673.D
Acq On : 28 Feb 2019 23:32
Operator : JU/SJ
Sample : K1643-03
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
FLOOR-SLAB

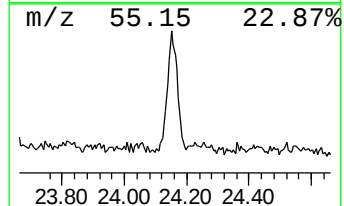
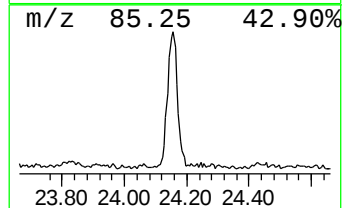
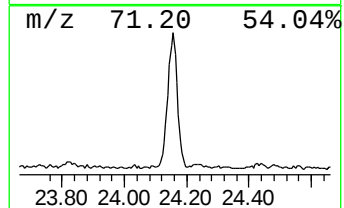
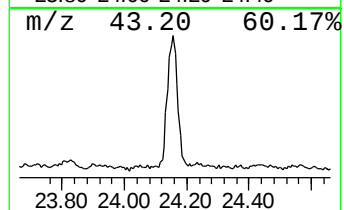
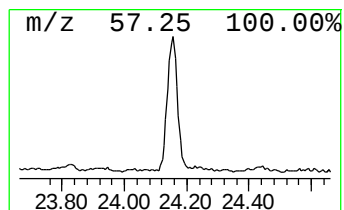
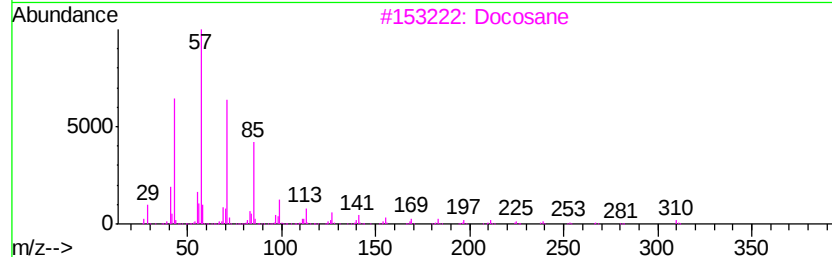
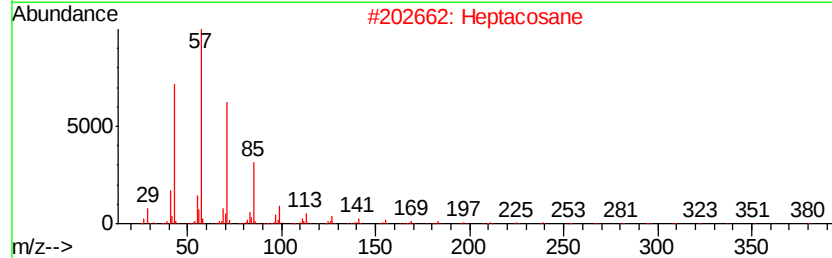
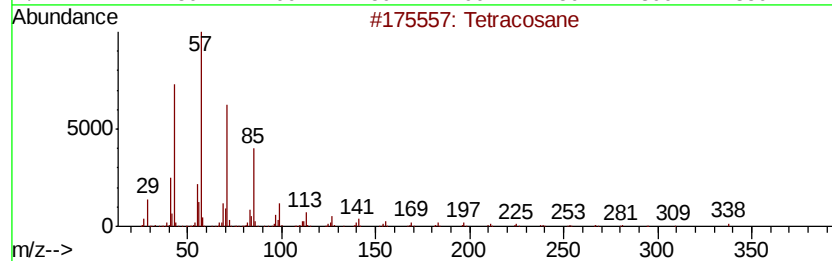
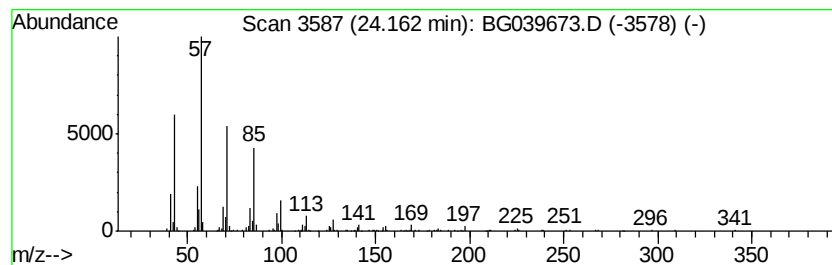
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 Tetracosane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.16	11.11 ng	503216	Perylene-d12	25.20

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetracosane	338	C24H50	000646-31-1	91
2		Heptacosane	380	C27H56	000593-49-7	91
3		Docosane	310	C22H46	000629-97-0	91
4		Eicosane	282	C20H42	000112-95-8	91
5		Heneicosane	296	C21H44	000629-94-7	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG022819\
Data File : BG039673.D
Acq On : 28 Feb 2019 23:32
Operator : JU/SJ
Sample : K1643-03
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
FLOOR-SLAB

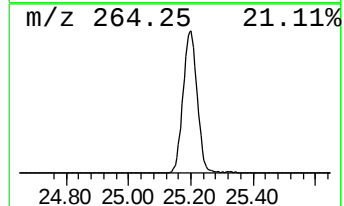
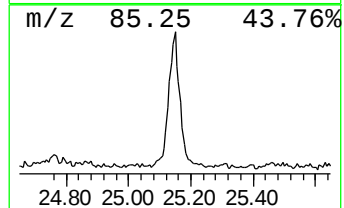
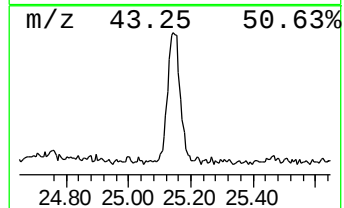
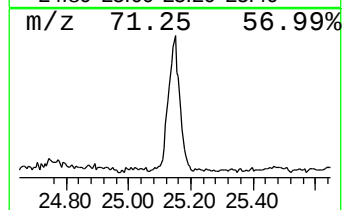
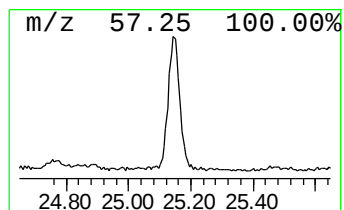
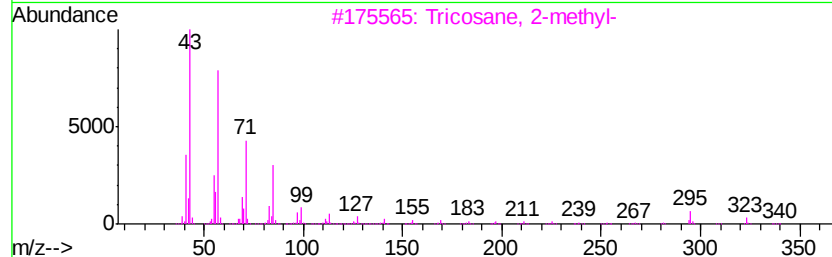
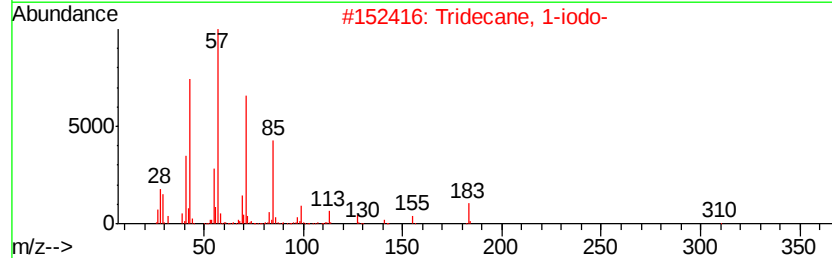
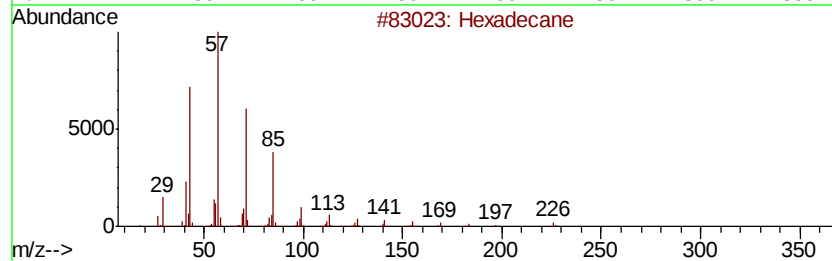
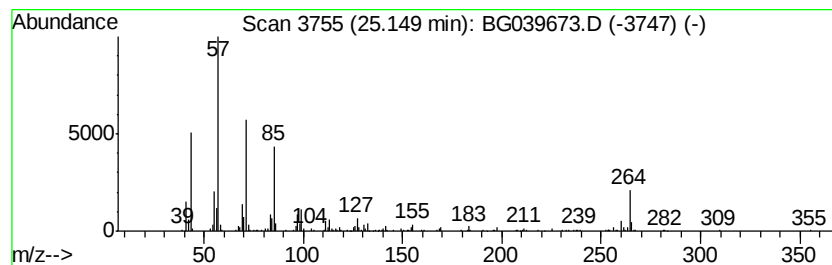
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 Hexadecane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.15	6.68 ng	302432	Perylene-d12	25.20

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexadecane	226	C16H34	000544-76-3	64
2			Tridecane, 1-iodo-	310	C13H27I	035599-77-0	64
3			Tricosane, 2-methyl-	338	C24H50	001928-30-9	58
4			Hentriacontane	437	C31H64	000630-04-6	53
5			1-Iodoundecane	282	C11H23I	004282-44-4	49



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG022819\
Data File : BG039673.D
Acq On : 28 Feb 2019 23:32
Operator : JU/SJ
Sample : K1643-03
Misc :
ALS Vial : 15 Sample Multiplier: 1

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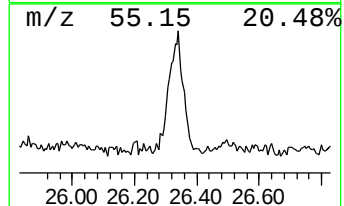
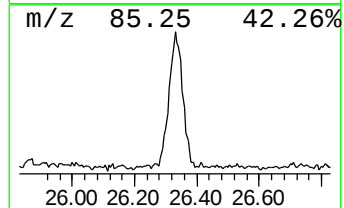
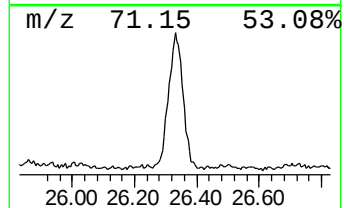
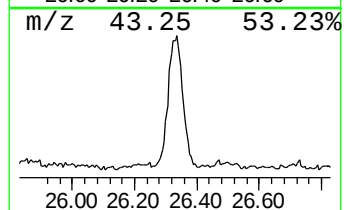
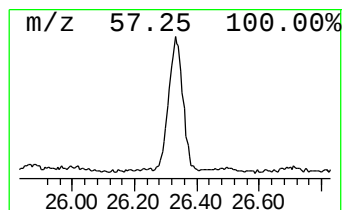
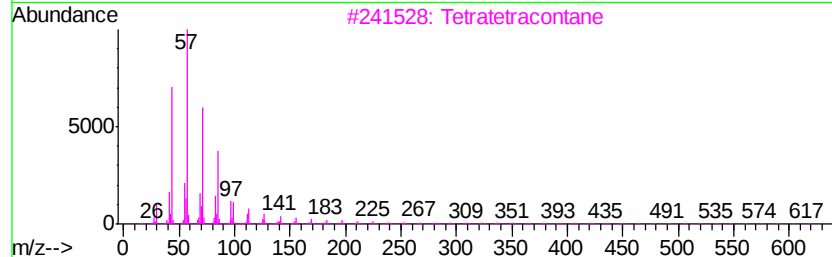
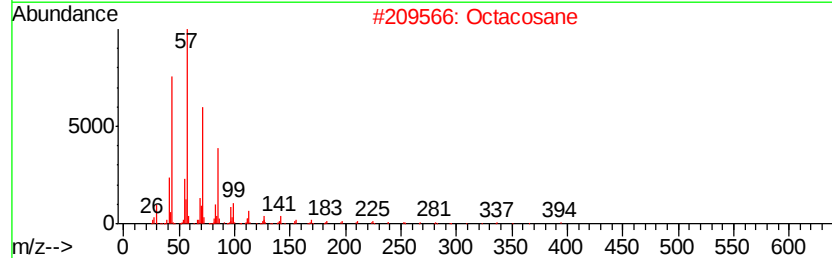
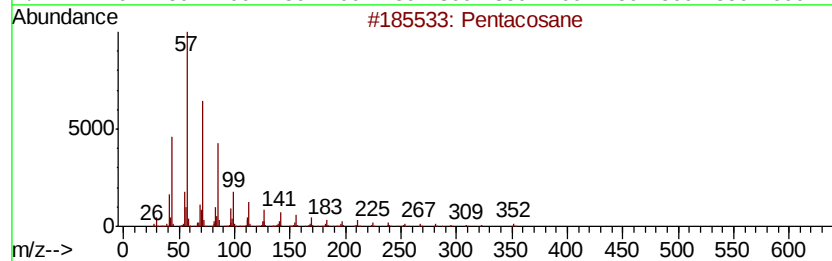
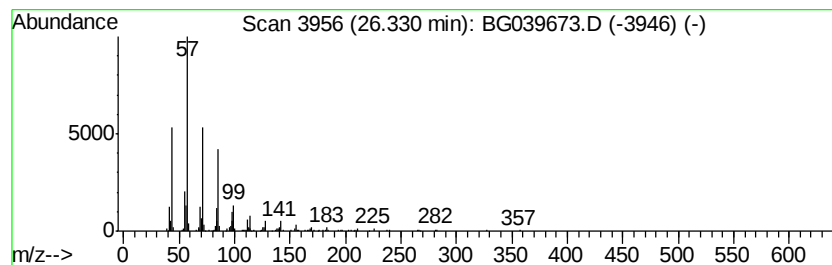
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 18 Pentacosane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.33	10.12 ng	458584	Perylene-d12	25.20

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentacosane	352	C25H52	000629-99-2	91
2		Octacosane	394	C28H58	000630-02-4	91
3		Tetratetracontane	619	C44H90	007098-22-8	91
4		Tetracosane, 11-decyl-	479	C34H70	055429-84-0	91
5		triacontane	422	C30H62	000638-68-6	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG022819\
Data File : BG039673.D
Acq On : 28 Feb 2019 23:32
Operator : JU/SJ
Sample : K1643-03
Misc :
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Instrument :
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ClientSampleId :
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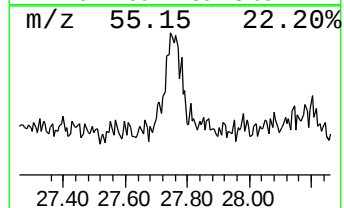
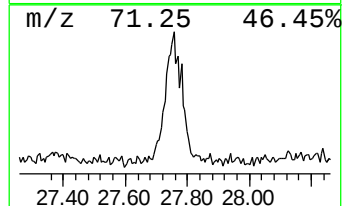
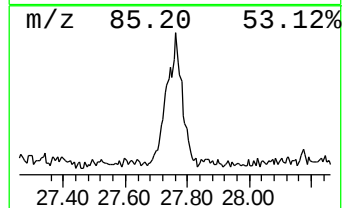
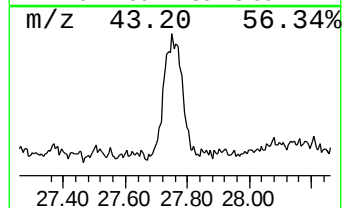
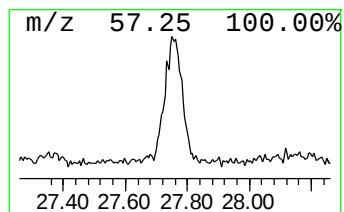
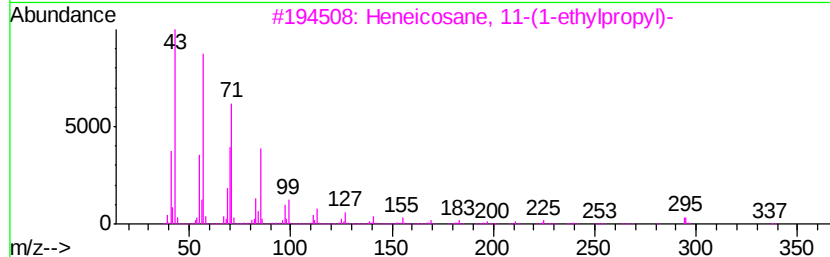
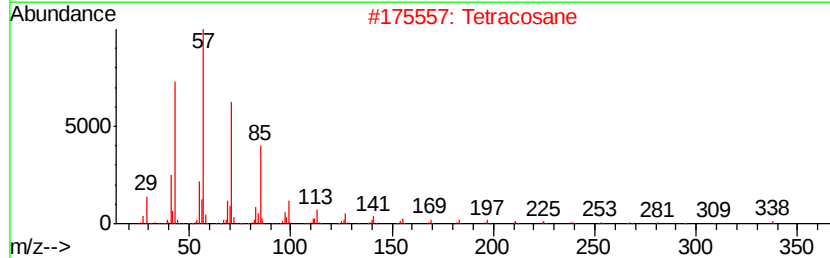
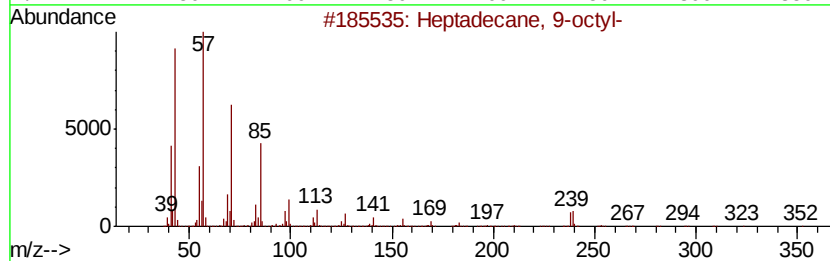
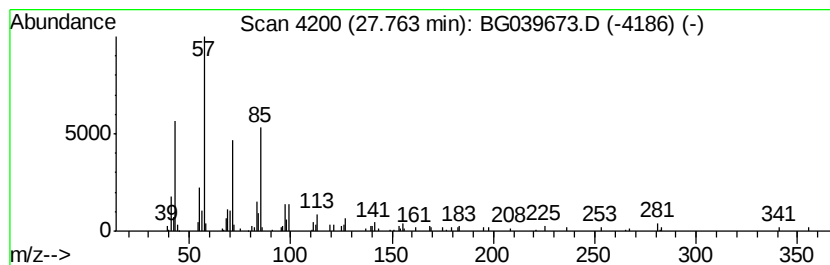
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 19 Heptadecane, 9-octyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
27.76	4.94 ng	223810	Perylene-d12	25.20

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	83
2		Tetracosane	338	C24H50	000646-31-1	83
3		Heneicosane, 11-(1-ethylpropyl)-	366	C26H54	055282-11-6	83
4		2-methyloctacosane	408	C29H60	1000376-72-8	83
5		Tetratetracontane	619	C44H90	007098-22-8	83



Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG022819\
 Data File : BG039673.D
 Acq On : 28 Feb 2019 23:32
 Operator : JU/SJ
 Sample : K1643-03
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
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 ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG012919.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
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3-Penten-2-one, 4...	4.64	11.8	ng	211930	1	8.17	358515	20.0
2-Pentanone, 4-hy...	5.25	22.7	ng	406602	1	8.17	358515	20.0
unknown7.70	7.70	103.6	ng	1856280	1	8.17	358515	20.0
unknown19.23	19.23	3.8	ng	187236	4	17.53	975808	20.0
Octacosane	19.29	2.9	ng	141201	4	17.53	975808	20.0
Eicosane	20.39	2.7	ng	129840	5	21.82	977047	20.0
Nonadecane, 9-met...	20.89	3.1	ng	149634	5	21.82	977047	20.0
Sulfurous acid, b...	21.41	4.8	ng	236567	5	21.82	977047	20.0
Nonadecane	21.97	4.4	ng	215733	5	21.82	977047	20.0
Heneicosane	22.60	6.5	ng	315064	5	21.82	977047	20.0
Heptadecane, 2,6,...	23.32	7.3	ng	354460	5	21.82	977047	20.0
Tetracosane	24.16	11.1	ng	503216	6	25.20	906099	20.0
Hexadecane	25.15	6.7	ng	302432	6	25.20	906099	20.0
Pentacosane	26.33	10.1	ng	458584	6	25.20	906099	20.0
Heptadecane, 9-oc...	27.76	4.9	ng	223810	6	25.20	906099	20.0