LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM051916\
Data File : BM005521.D
Acq On : 20 May 2016 05:11
Operator : UM/SJ
Sample : H2841-16
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : D9WT1

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 1 % 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

Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM051816.M
Title : SVOA CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	7.004	742	751	760	rBV2	148480	280632	3.62%	0.475%
2	7.169	773	779	789	rBV2	126911	229902	2.96%	0.389%
3	7.369	807	813	820	rBV	159530	270013	3.48%	0.457%
4	7.839	885	893	902	rBV	582875	978782	12.62%	1.656%
5	8.539	1002	1012	1022	rBV	226918	382149	4.93%	0.646%
6	8.745	1039	1047	1057	rBV3	30687	100144	1.29%	0.169%
7	8.992	1083	1089	1098	rBV	160331	273250	3.52%	0.462%
8	9.710	1205	1211	1219	rBV	144193	251062	3.24%	0.425%
9	10.251	1296	1303	1311	rVB	247024	425764	5.49%	0.720%
10	10.627	1358	1367	1374	rBV	959000	1647986	21.25%	2.788%
11	10.763	1386	1390	1395	rBV	54834	104272	1.34%	0.176%
12	10.810	1395	1398	1406	rVB2	49138	79135	1.02%	0.134%
13	11.904	1577	1584	1589	rVB	114607	172969	2.23%	0.293%
14	12.216	1631	1637	1644	rBV2	239728	414234	5.34%	0.701%
15	12.286	1644	1649	1655	rVB	71336	115980	1.50%	0.196%
16	13.639	1873	1879	1886	rVB5	83673	155950	2.01%	0.264%
17	13.792	1898	1905	1910	rBV2	50670	87753	1.13%	0.148%
18	13.880	1915	1920	1925	rVV	504617	899473	11.60%	1.522%
19	13.921	1925	1927	1936	rVB	121626	172366	2.22%	0.292%
20	14.162	1961	1968	1981	rVB	561753	891011	11.49%	1.507%
21	14.468	2012	2020	2028	rBV2	1589670	2527697	32.59%	4.276%
22	14.662	2047	2053	2064	rBV2	170997	307871	3.97%	0.521%
23	14.874	2084	2089	2095	rBV5	53208	104440	1.35%	0.177%
24	15.315	2161	2164	2169	rVB	65794	86887	1.12%	0.147%
25	15.457	2183	2188	2194	rBV2	864859	1297273	16.73%	2.195%
26	15.574	2202	2208	2215	rBV	175539	279673	3.61%	0.473%
27	15.721	2225	2233	2240	rBV4	76107	161979	2.09%	0.274%
28	16.180	2307	2311	2313	rBV2	77508	113050	1.46%	0.191%
29	16.204	2313	2315	2320	rVB2	104083	129320	1.67%	0.219%
30	16.356	2337	2341	2347	rBV3	63532	94368	1.22%	0.160%
31	16.415	2347	2351	2355	rBV2	57324	78443	1.01%	0.133%
32	16.562	2373	2376	2380	rVB	117324	134880	1.74%	0.228%
33	16.621	2380	2386	2393	rBV4	410926	817347	10.54%	1.383%
34	16.703	2396	2400	2404	rVV3	114839	184039	2.37%	0.311%

LSC Area Percent Report

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Data File : BM005521.D
Acq On : 20 May 2016 05:11
Operator : UM/SJ
Sample : H2841-16
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : D9WT1

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 1 % 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

Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM051816.M
Title : SVOA CALIBRATION

35	16.756	2405	2409	2415	rVB	107967	149024	1.92%	0.252%
36	17.109	2466	2469	2475	rBV8	52883	103245	1.33%	0.175%
37	17.209	2478	2486	2497	rVV2	2327943	3724730	48.03%	6.301%
38	17.309	2497	2503	2509	rBV2	959355	1437566	18.54%	2.432%
39	17.592	2547	2551	2559	rBV2	380574	566144	7.30%	0.958%
40	18.109	2635	2639	2646	rBV	627787	923980	11.91%	1.563%
41	18.627	2723	2727	2733	rVB	575814	801943	10.34%	1.357%
42	18.962	2780	2784	2788	rBV	970258	1241357	16.01%	2.100%
43	19.297	2838	2841	2849	rVB	423375	581141	7.49%	0.983%
44	19.386	2854	2856	2866	rVB2	460606	728060	9.39%	1.232%
45	19.603	2889	2893	2898	rBV	1139802	1494496	19.27%	2.528%
46	21.391	3193	3197	3204	rVB	3576799	4638875	59.82%	7.848%
47	21.597	3229	3232	3236	rVB	1309929	1570982	20.26%	2.658%
48	21.709	3248	3251	3256	rVB	2228827	2897929	37.37%	4.902%
49	23.697	3585	3589	3609	rVB	2697582	6061288	78.16%	10.254%
50	24.327	3691	3696	3709	rVB2	3491835	7755185	100.00%	13.120%
51	24.562	3730	3736	3742	rVB	2273378	4665181	60.16%	7.892%
52	26.073	3987	3993	4006	rVB3	1852251	5520041	71.18%	9.338%

Sum of corrected areas: 59111261

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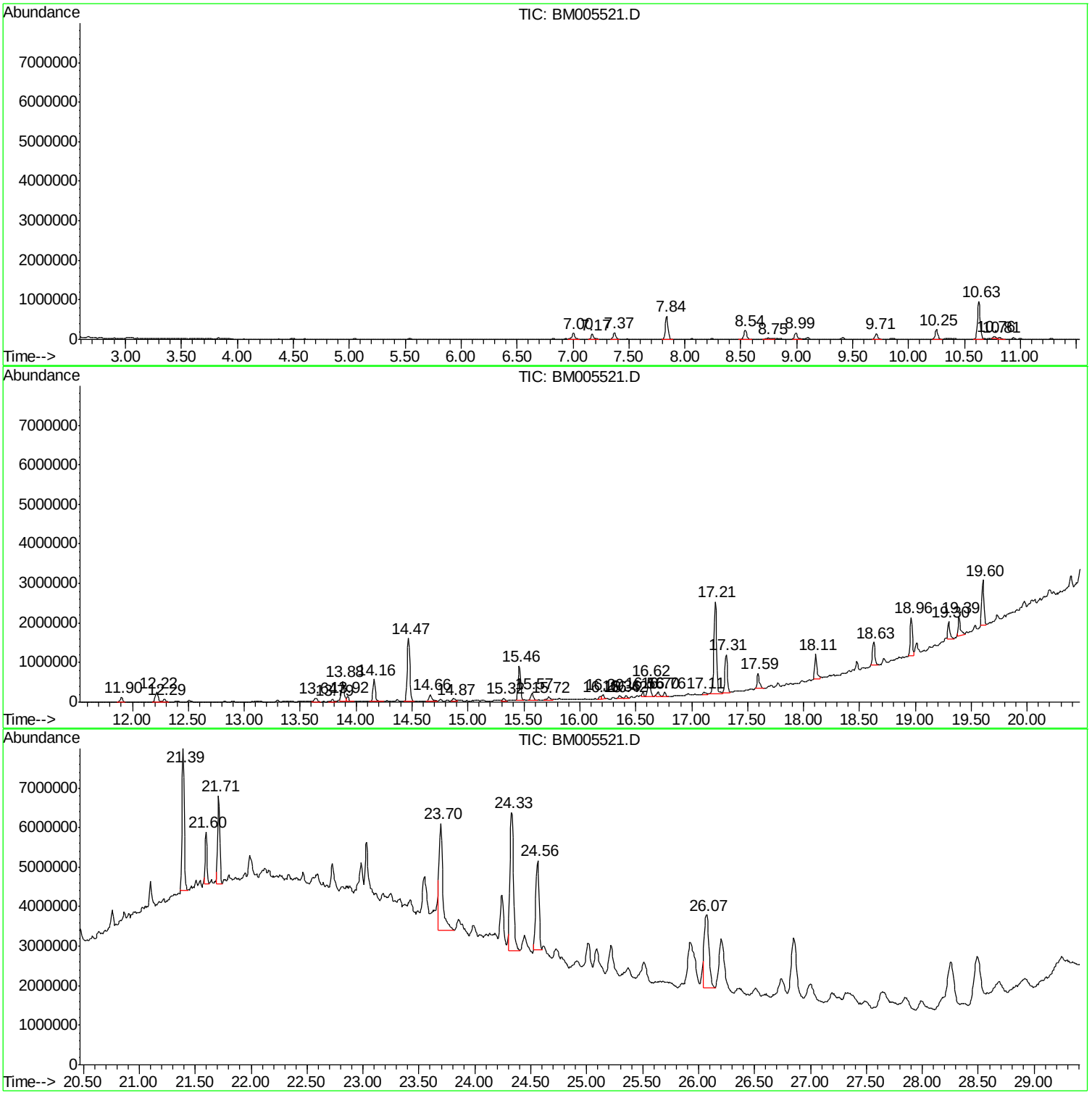
Acq On : 20 May 2016 05:11
Operator : UM/SJ
Sample : H2841-16
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
D9WT1

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM051816.M
Quant Title : SVOA CALIBRATION

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



Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM051916\
Data File : BM005521.D
Acq On : 20 May 2016 05:11
Operator : UM/SJ
Sample : H2841-16
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : D9WT1

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM051816.M
Quant Title : SVOA CALIBRATION

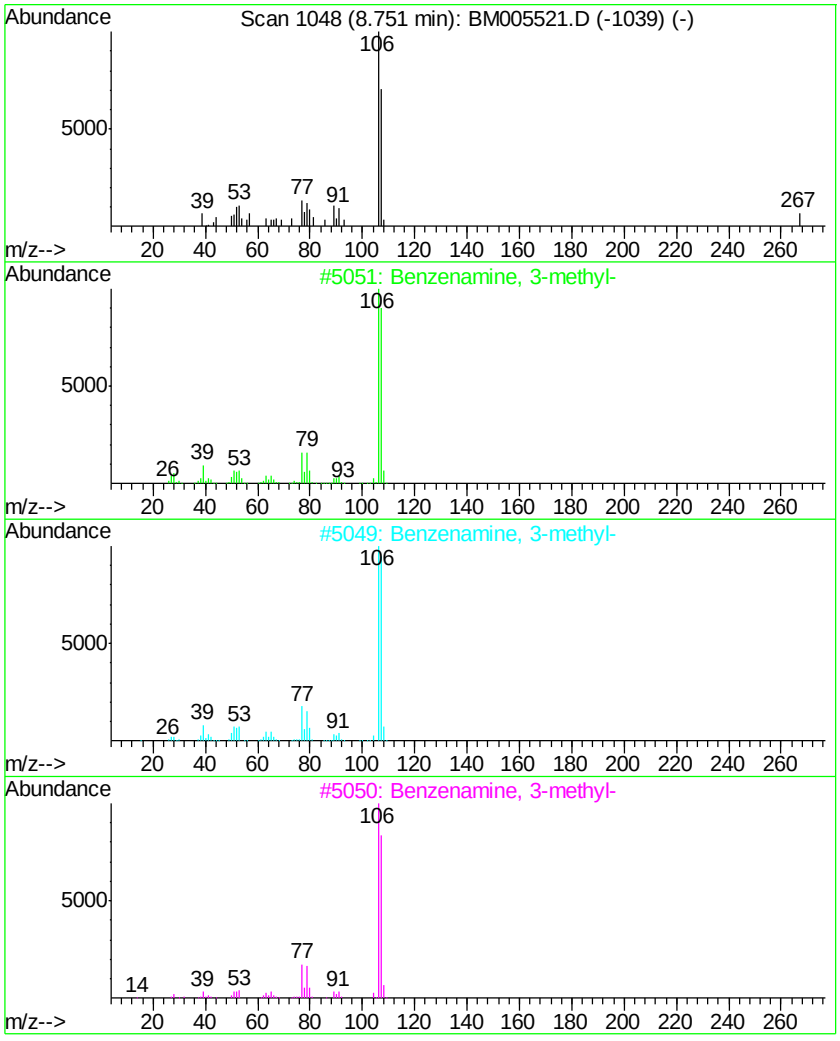
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

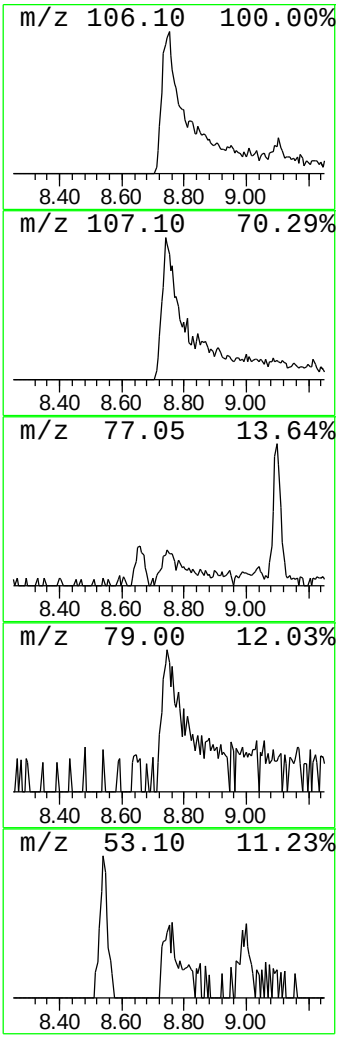
****

Peak Number 1 Benzenamine, 3-methyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.75	2.05 ng/ul	100144	1,4-Dichlorobenzene-d4	7.84

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzenamine, 3-methyl-	107	C7H9N	000108-44-1	87
2	Benzenamine, 3-methyl-	107	C7H9N	000108-44-1	83
3	Benzenamine, 3-methyl-	107	C7H9N	000108-44-1	80
4	o-Toluidine	107	C7H9N	000095-53-4	74
5	Urea, (3-methylphenyl)-	150	C8H10N2O	000063-99-0	72





Library Search Compound Report

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Data File : BM005521.D
Acq On : 20 May 2016 05:11
Operator : UM/SJ
Sample : H2841-16
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : D9WT1

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM051816.M
Quant Title : SVOA CALIBRATION

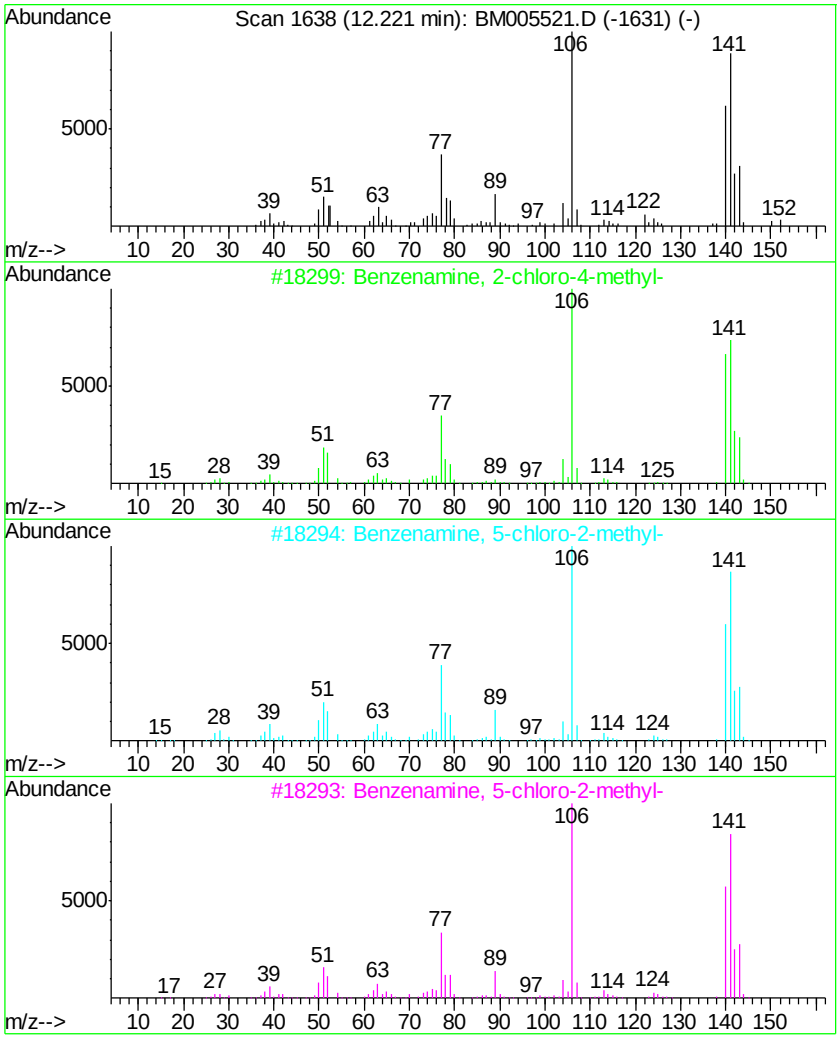
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

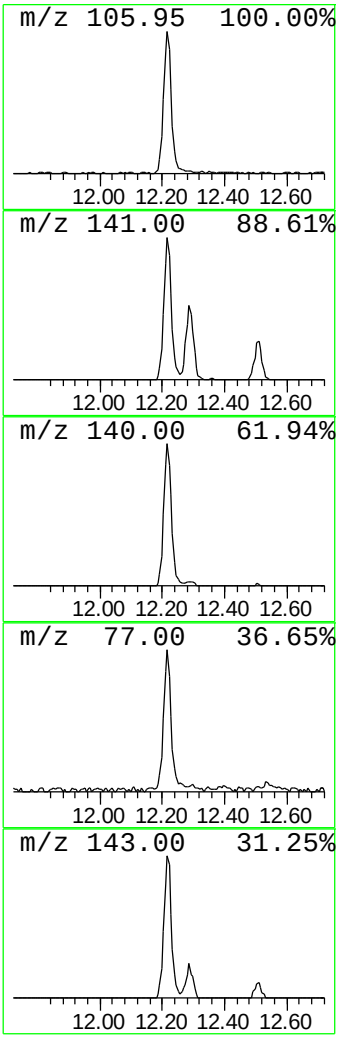
****

Peak Number 3 Benzenamine, 2-chloro-4-met... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.22	5.03 ng/ul	414234	Naphthalene-d8	10.63

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzenamine, 2-chloro-4-methyl-	141	C7H8ClN	000615-65-6	98
2	Benzenamine, 5-chloro-2-methyl-	141	C7H8ClN	000095-79-4	98
3	Benzenamine, 5-chloro-2-methyl-	141	C7H8ClN	000095-79-4	98
4	Benzenamine, 2-chloro-4-methyl-	141	C7H8ClN	000615-65-6	97
5	2-Chloro-5-methylaniline	141	C7H8ClN	000095-81-8	95





Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM051916\
Data File : BM005521.D
Acq On : 20 May 2016 05:11
Operator : UM/SJ
Sample : H2841-16
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : D9WT1

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM051816.M
Quant Title : SVOA CALIBRATION

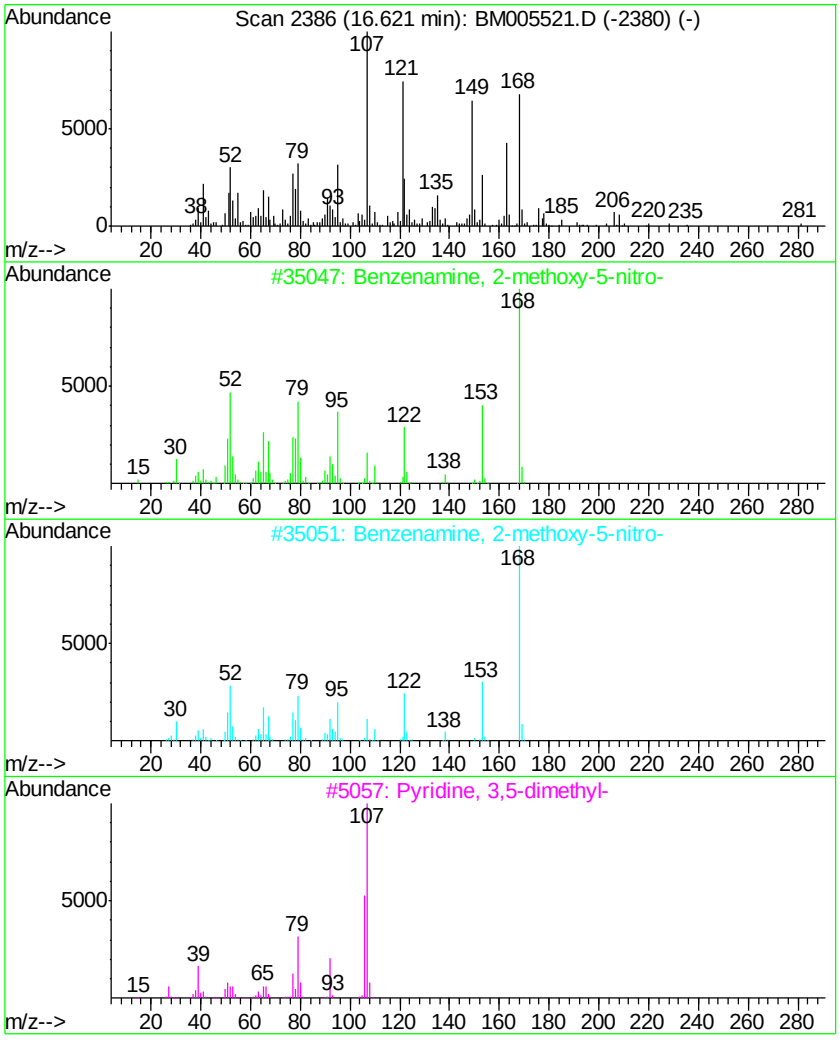
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

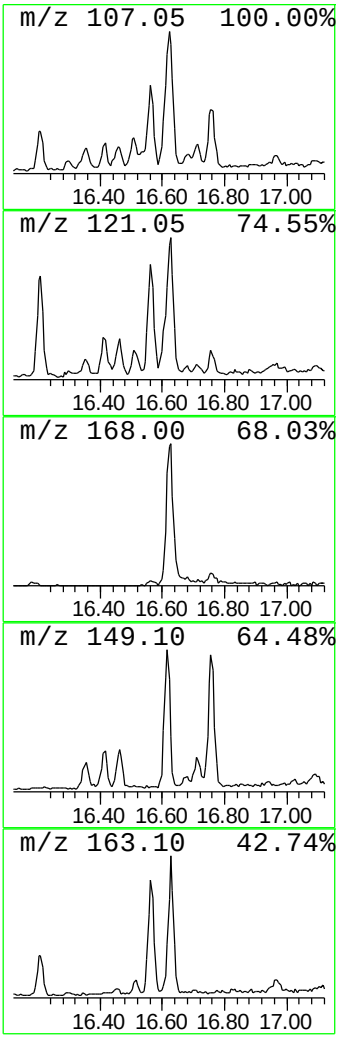
****

Peak Number 4 Benzenamine, 2-methoxy-5-ni... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.62	4.39 ng/ul	817347	Phenanthrene-d10	17.21

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzenamine, 2-methoxy-5-nitro-	168	C7H8N2O3	000099-59-2	94
2	Benzenamine, 2-methoxy-5-nitro-	168	C7H8N2O3	000099-59-2	42
3	Pyridine, 3,5-dimethyl-	107	C7H9N	000591-22-0	38
4	Benzenamine, 4-methoxy-2-nitro-	168	C7H8N2O3	000096-96-8	38
5	Benzenamine, 4-methoxy-2-nitro-	168	C7H8N2O3	000096-96-8	30





Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM051916\
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Operator : UM/SJ
Sample : H2841-16
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : D9WT1

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM051816.M
Quant Title : SVOA CALIBRATION

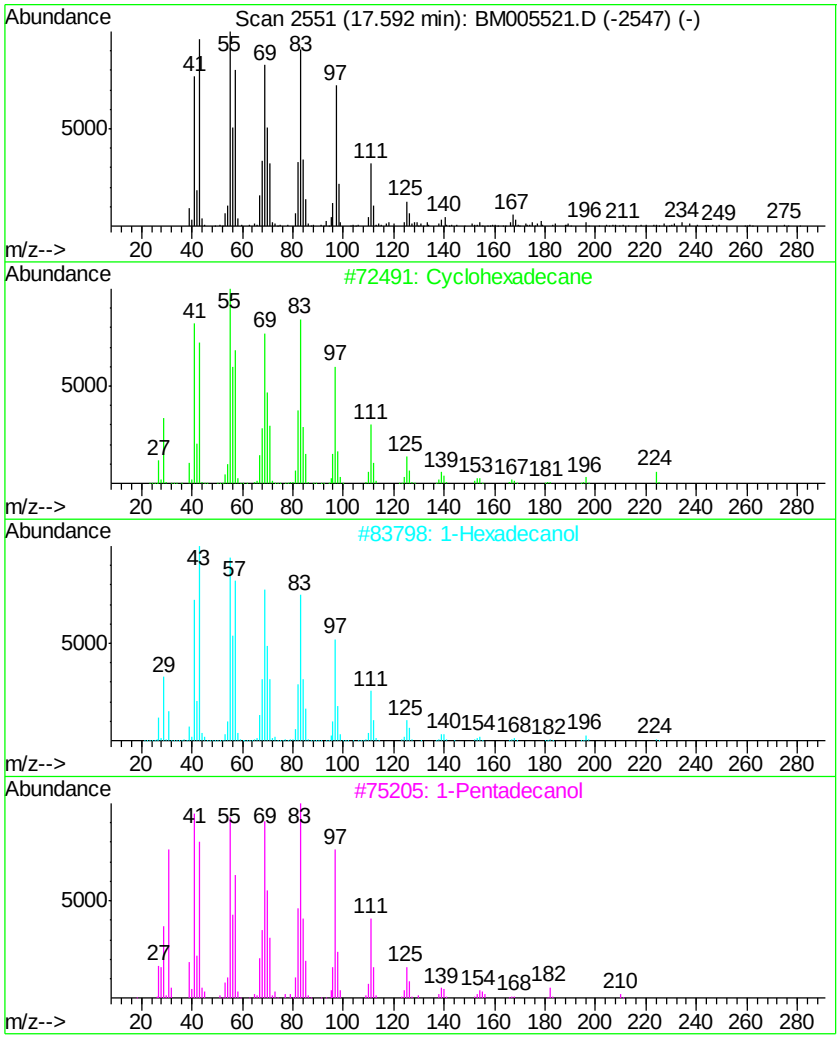
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

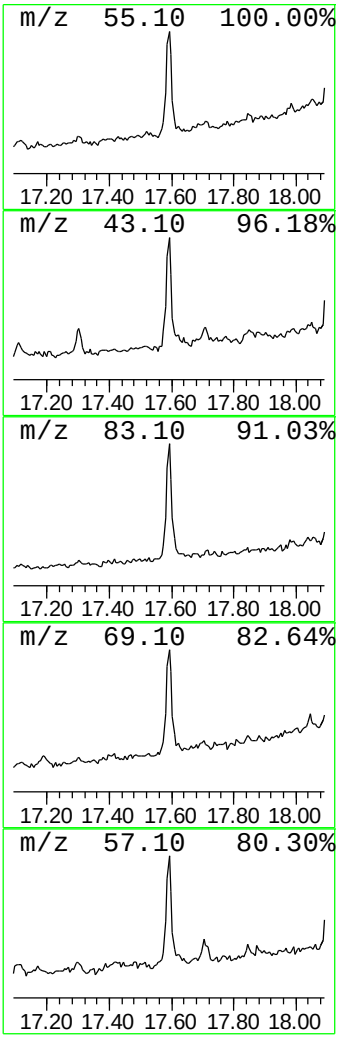
****

Peak Number 5 (DEL) Alkane: Cyclic-17.59 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.59	3.04 ng/ul	566144	Phenanthrene-d10	17.21

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexadecane	224	C16H32	000295-65-8	96
2	1-Hexadecanol	242	C16H34O	036653-82-4	94
3	1-Pentadecanol	228	C15H32O	000629-76-5	91
4	1-Tetradecanol	214	C14H30O	000112-72-1	91
5	7-Hexadecene, (Z)-	224	C16H32	035507-09-6	91





Data Path : Z:\HPCHEM1\BNA_M\DATA\BM051916\
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Operator : UM/SJ
Sample : H2841-16
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : D9WT1

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM051816.M
Quant Title : SVOA CALIBRATION

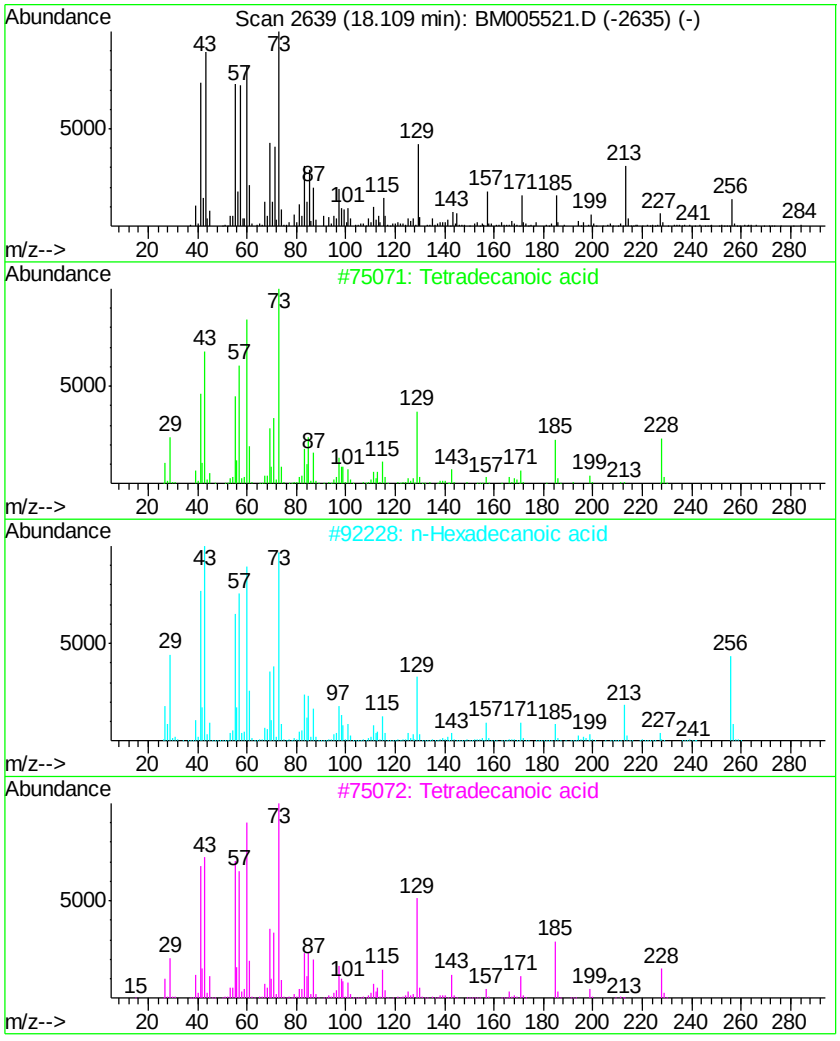
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

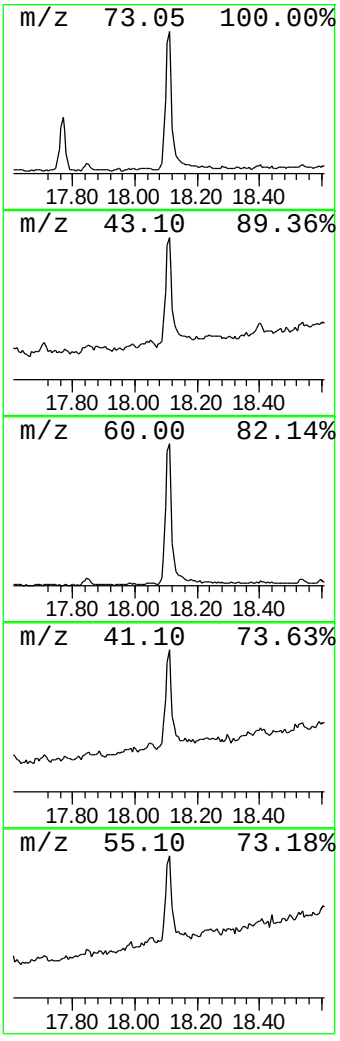
****

Peak Number 6 Tetradecanoic acid Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.11	4.96 ng/ul	923980	Phenanthrene-d10	17.21

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetradecanoic acid	228	C14H28O2	000544-63-8	95
2	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	95
3	Tetradecanoic acid	228	C14H28O2	000544-63-8	93
4	Undecanoic acid	186	C11H22O2	000112-37-8	83
5	Pentadecanoic acid	242	C15H30O2	001002-84-2	72





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Sample : H2841-16
Misc :
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Instrument : BNA_M
ClientSampleId : D9WT1

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM051816.M
Quant Title : SVOA CALIBRATION

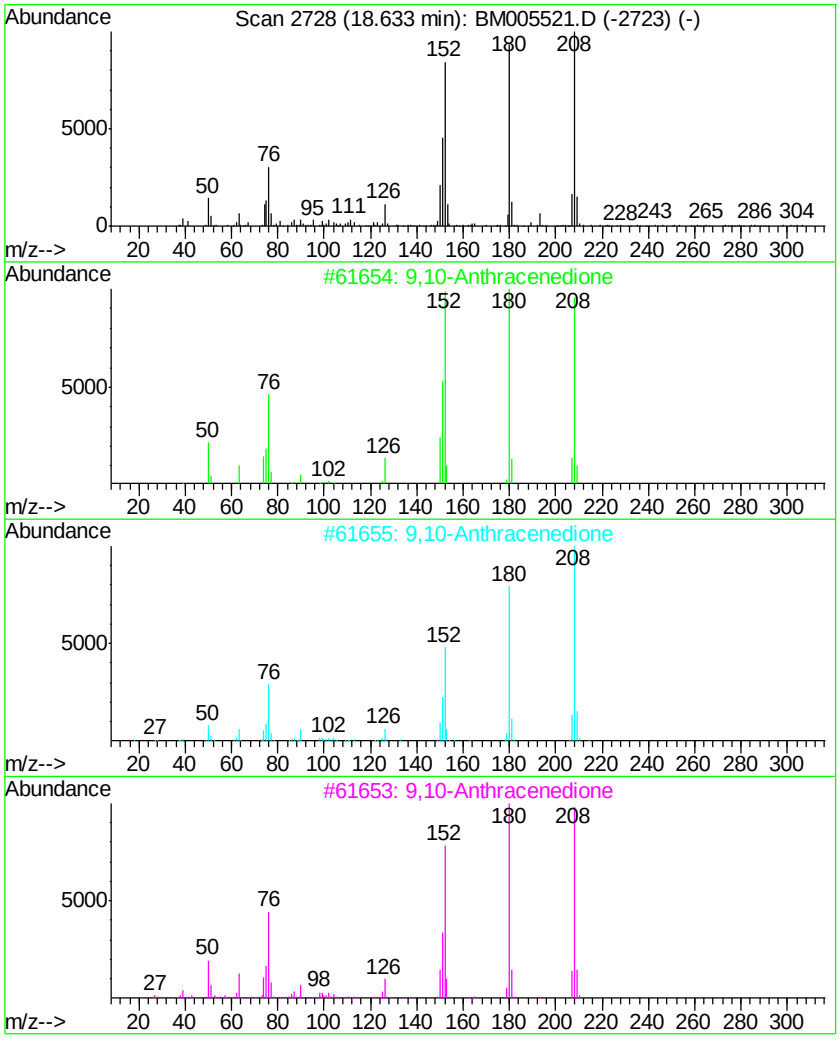
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

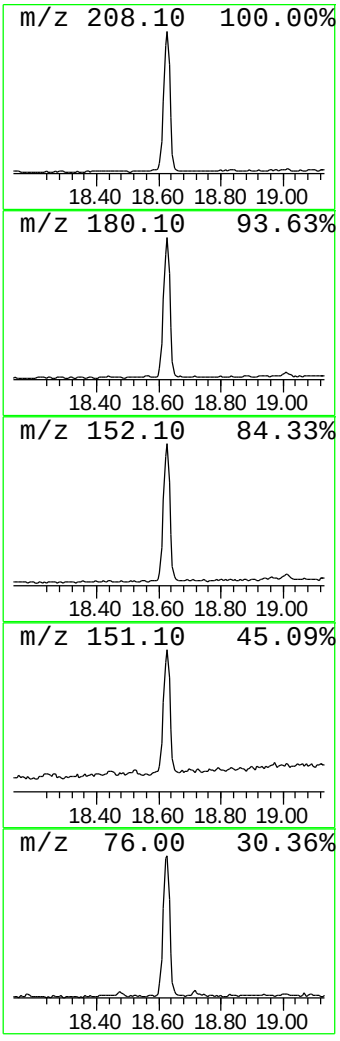
****

Peak Number 7 9,10-Anthracenedione Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.63	4.31 ng/ul	801943	Phenanthrene-d10	17.21

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9,10-Anthracenedione	208	C14H8O2	000084-65-1	99
2	9,10-Anthracenedione	208	C14H8O2	000084-65-1	95
3	9,10-Anthracenedione	208	C14H8O2	000084-65-1	94
4	Benzo[c]cinnoline	180	C12H8N2	000230-17-1	93
5	Benzo[c]cinnoline	180	C12H8N2	000230-17-1	86





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Sample : H2841-16
Misc :
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Instrument : BNA_M
ClientSampleId : D9WT1

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM051816.M
Quant Title : SVOA CALIBRATION

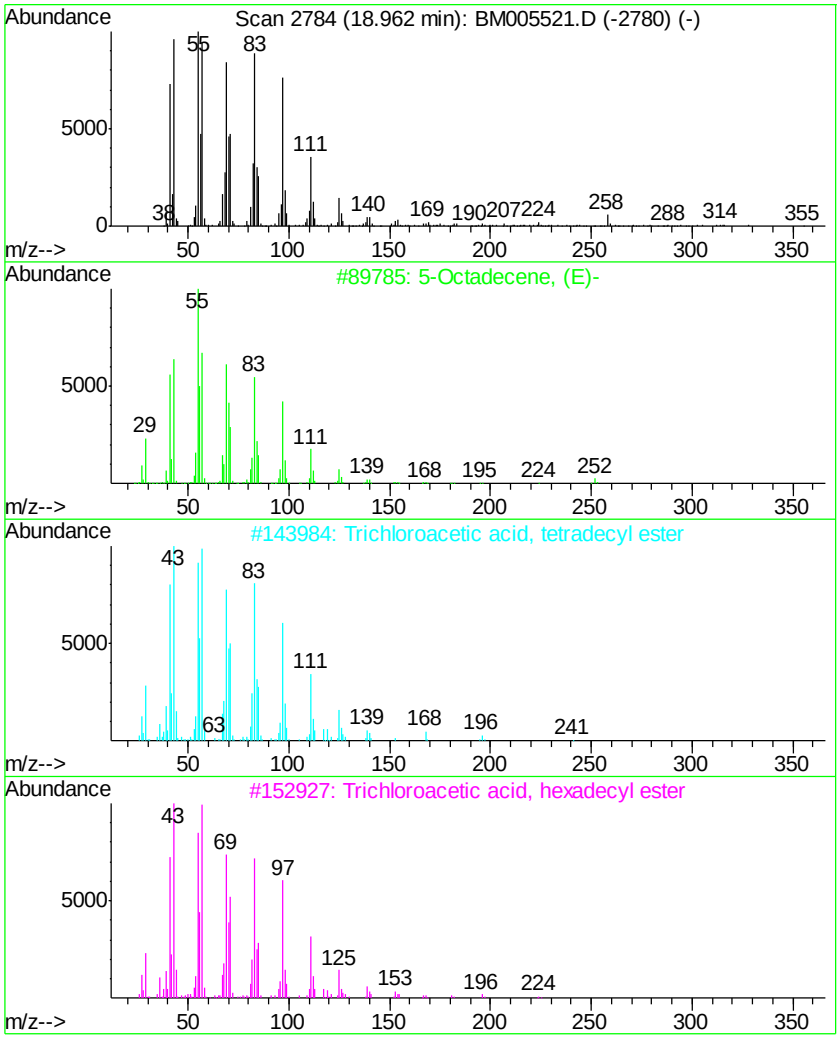
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

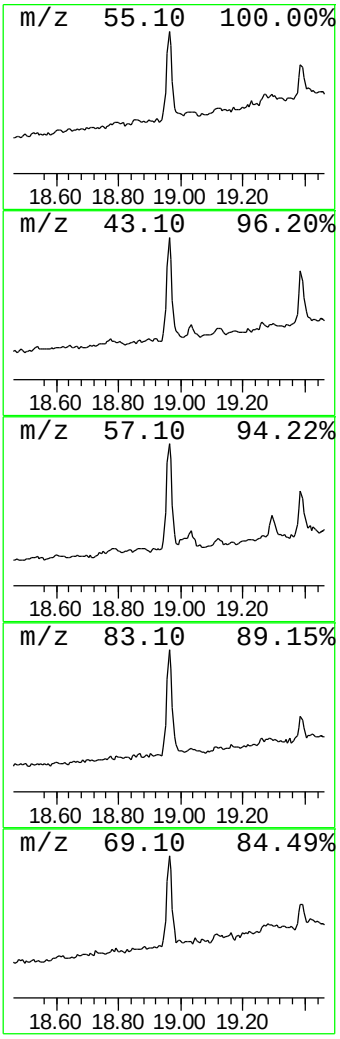
****

Peak Number 8 (DEL) Alkane: Cyclic-18.96 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.96	6.67 ng/ul	1241360	Phenanthrene-d10	17.21

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	5-Octadecene, (E)-	252	C18H36	007206-21-5	98
2	Trichloroacetic acid, tetradecyl...	358	C16H29Cl3O2	074339-52-9	94
3	Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	94
4	Cyclohexadecane	224	C16H32	000295-65-8	93
5	1-Octadecene	252	C18H36	000112-88-9	93





Data Path : Z:\HPCHEM1\BNA_M\DATA\BM051916\
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Operator : UM/SJ
Sample : H2841-16
Misc :
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Instrument : BNA_M
ClientSampleId : D9WT1

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM051816.M
Quant Title : SVOA CALIBRATION

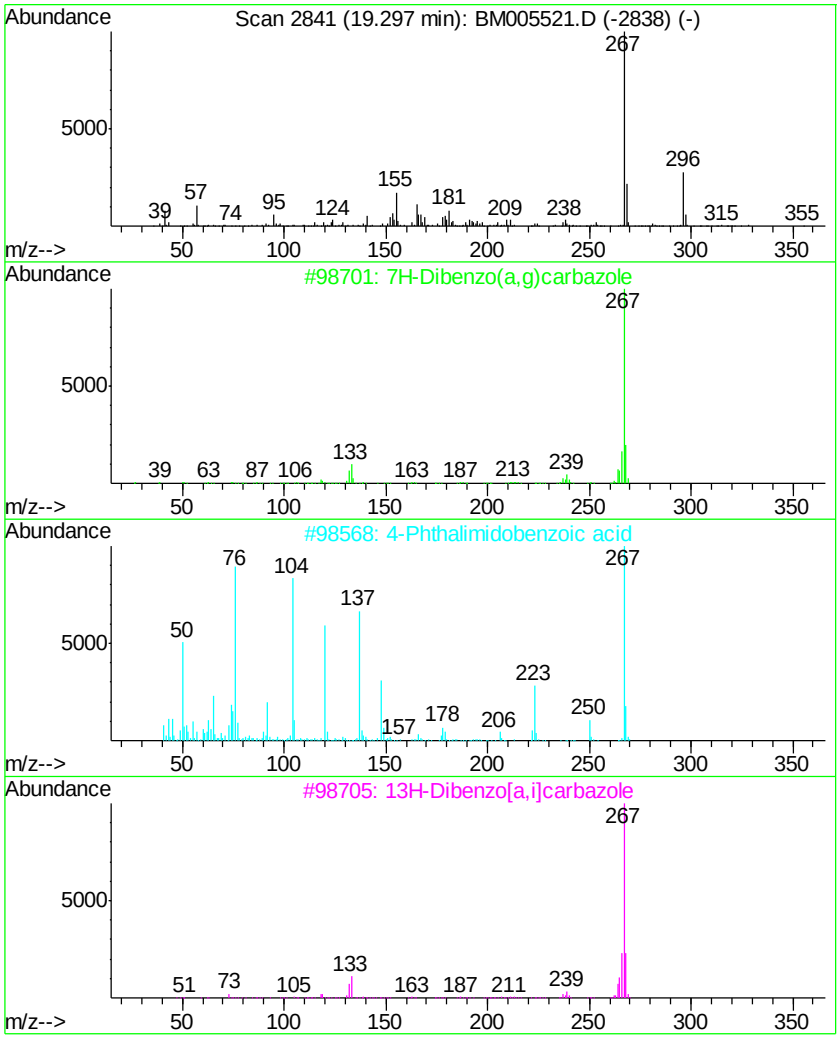
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

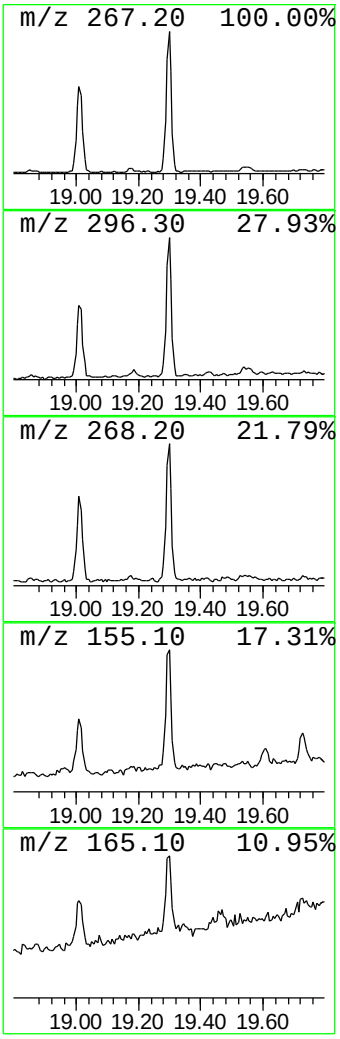
****

Peak Number 9 7H-Dibenzo(a,g)carbazole Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.30	3.12 ng/ul	581141	Phenanthrene-d10	17.21

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	7H-Dibenzo(a,g)carbazole	267	C20H13N	000207-84-1	52
2	4-Phthalimidobenzoic acid	267	C15H9NO4	005383-82-4	52
3	13H-Dibenzo[a,i]carbazole	267	C20H13N	000239-64-5	50
4	13H-Dibenzo[a,i]carbazole	267	C20H13N	000239-64-5	50
5	5H-Naphtho[2,3-b]carbazole	267	C20H13N	000248-96-4	50





Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM051916\
Data File : BM005521.D
Acq On : 20 May 2016 05:11
Operator : UM/SJ
Sample : H2841-16
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : D9WT1

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM051816.M
Quant Title : SVOA CALIBRATION

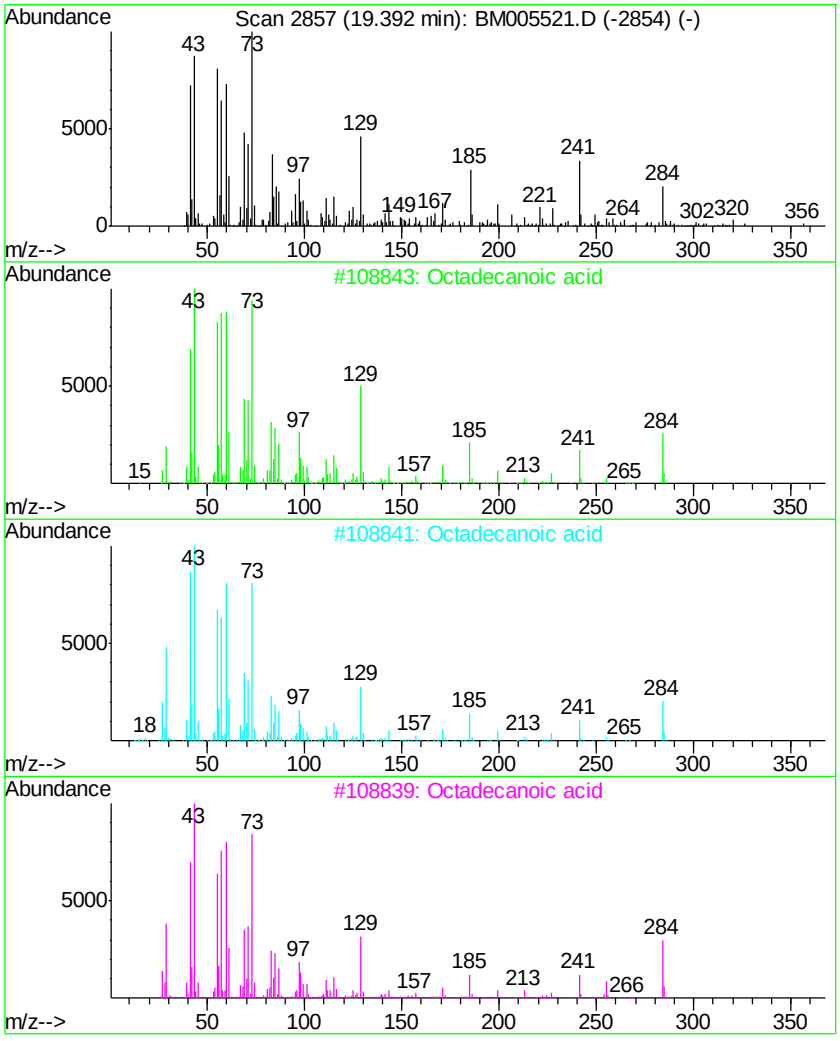
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

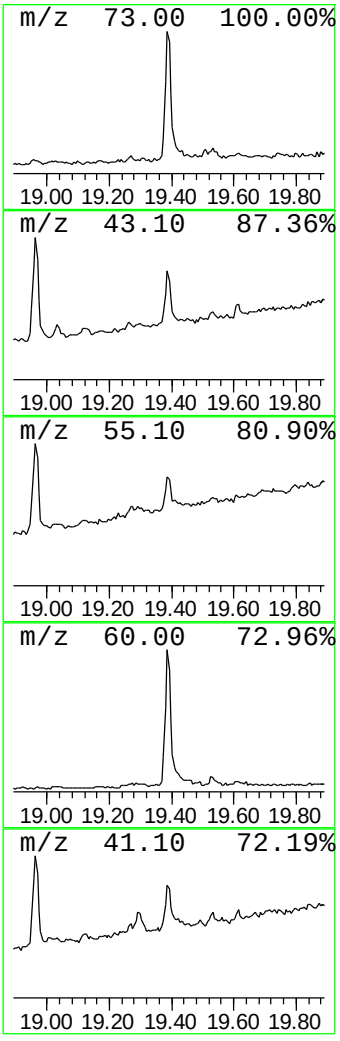
****

Peak Number 10 Octadecanoic acid Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.39	3.14 ng/ul	728060	Chrysene-d12	21.39

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecanoic acid	284	C18H36O2	000057-11-4	97
2	Octadecanoic acid	284	C18H36O2	000057-11-4	96
3	Octadecanoic acid	284	C18H36O2	000057-11-4	93
4	Octadecanoic acid	284	C18H36O2	000057-11-4	91
5	Pentadecanoic acid	242	C15H30O2	001002-84-2	89





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Operator : UM/SJ
Sample : H2841-16
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : D9WT1

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM051816.M
Quant Title : SVOA CALIBRATION

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

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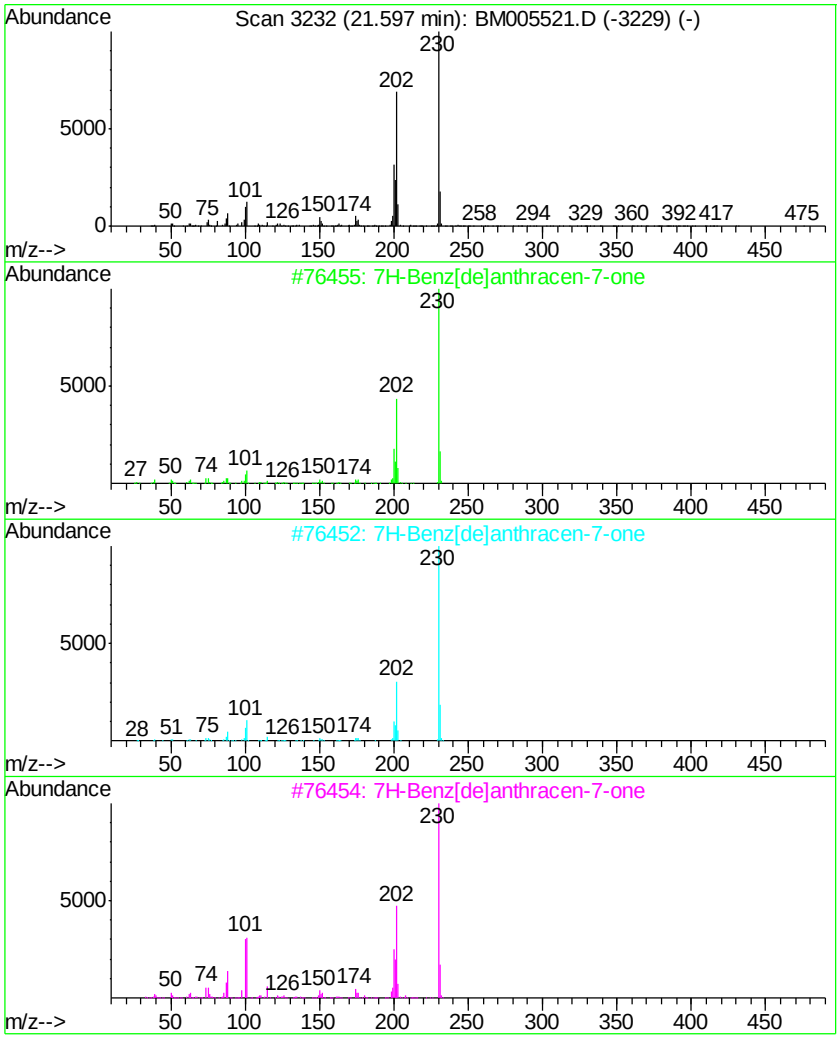
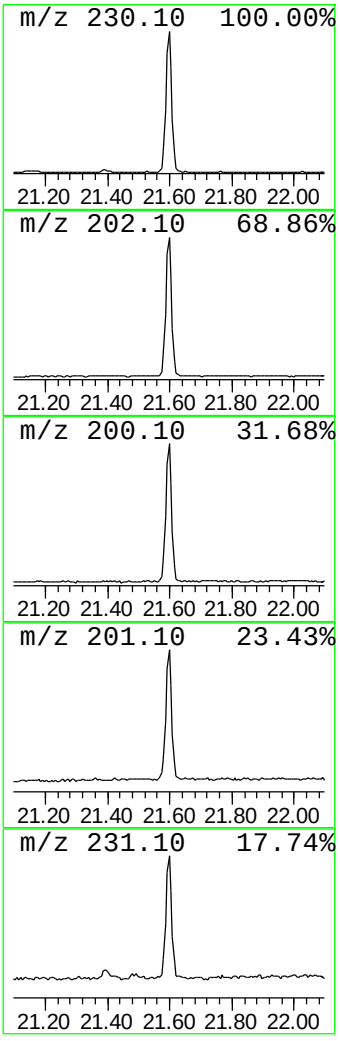
Peak Number 11 7H-Benz[de]anthracen-7-one Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.60	6.77 ng/ul	1570980	Chrysene-d12	21.39

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	7H-Benz[de]anthracen-7-one	230	C17H10O	000082-05-3	97
2	7H-Benz[de]anthracen-7-one	230	C17H10O	000082-05-3	93
3	7H-Benz[de]anthracen-7-one	230	C17H10O	000082-05-3	90
4	7H-Benz[de]anthracen-7-one	230	C17H10O	000082-05-3	86
5	Fluoranthene	202	C16H10	000206-44-0	83





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Data File : BM005521.D
Acq On : 20 May 2016 05:11
Operator : UM/SJ
Sample : H2841-16
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : D9WT1

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM051816.M
Quant Title : SVOA CALIBRATION

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

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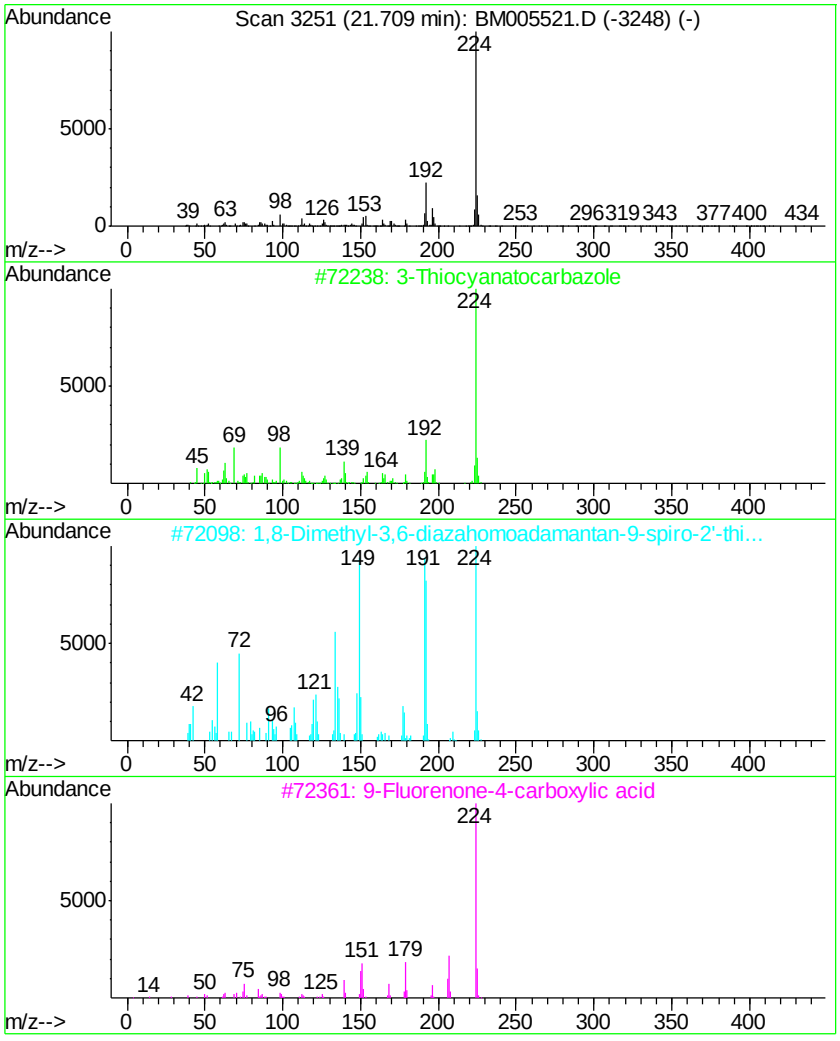
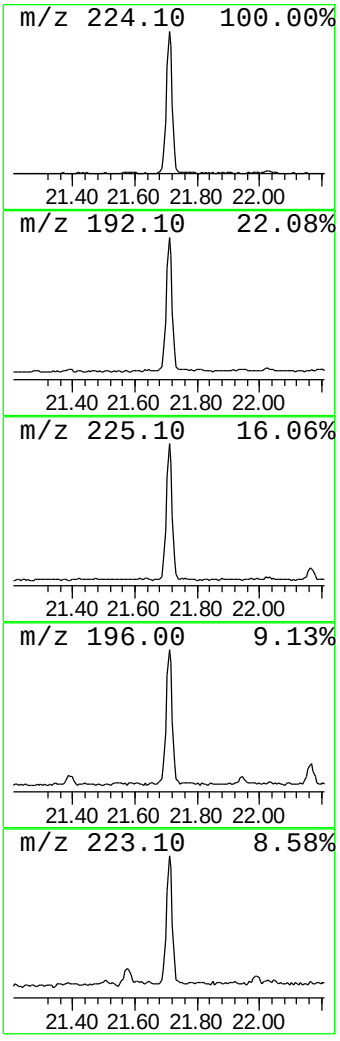
Peak Number 12 3-Thiocyanatocarbazole Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.71	12.49 ng/ul	2897930	Chrysene-d12	21.39

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Thiocyanatocarbazole	224	C13H8N2S	040604-35-1	90
2	1,8-Dimethyl-3,6-diazahomoadaman...	224	C12H20N2S	1000216-30-4	64
3	9-Fluorenone-4-carboxylic acid	224	C14H8O3	006223-83-2	53
4	Dibenzo[d,f]-1,3,2-dioxaborepine...	224	C14H13BO2	1000160-51-1	53
5	2-Propen-1-one, 1-(2-hydroxyphen...	224	C15H12O2	001214-47-7	50





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Acq On : 20 May 2016 05:11
Operator : UM/SJ
Sample : H2841-16
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : D9WT1

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM051816.M
Quant Title : SVOA CALIBRATION

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

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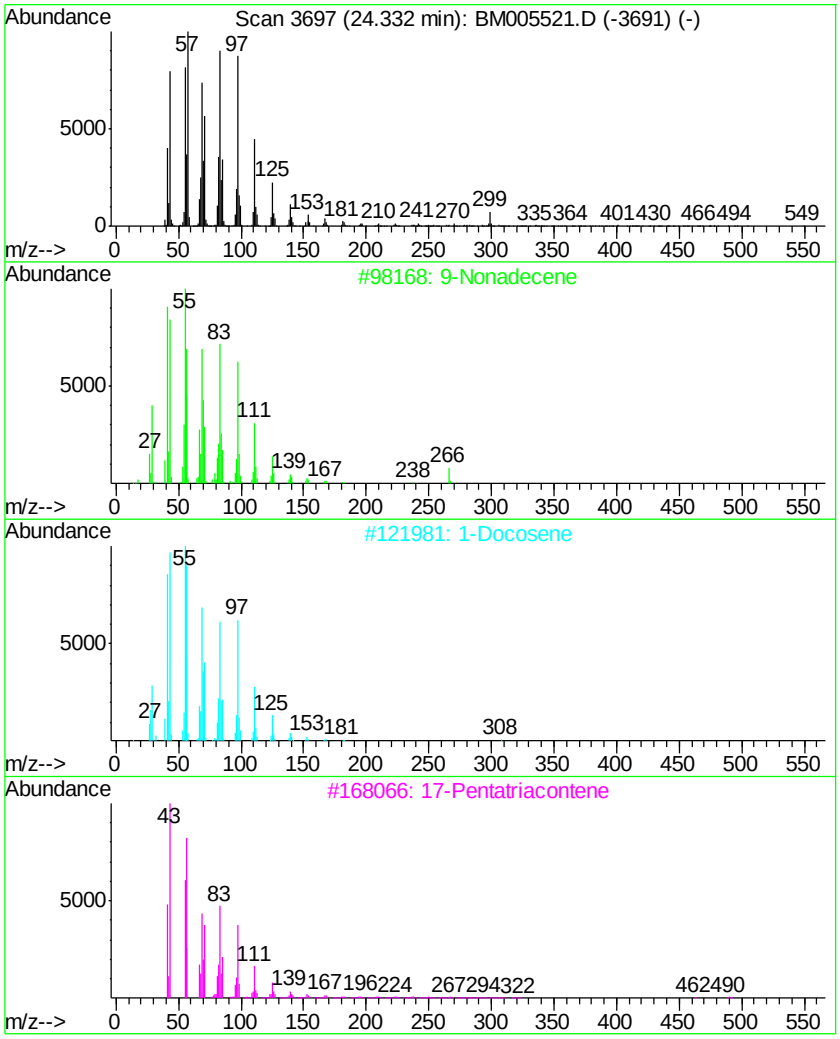
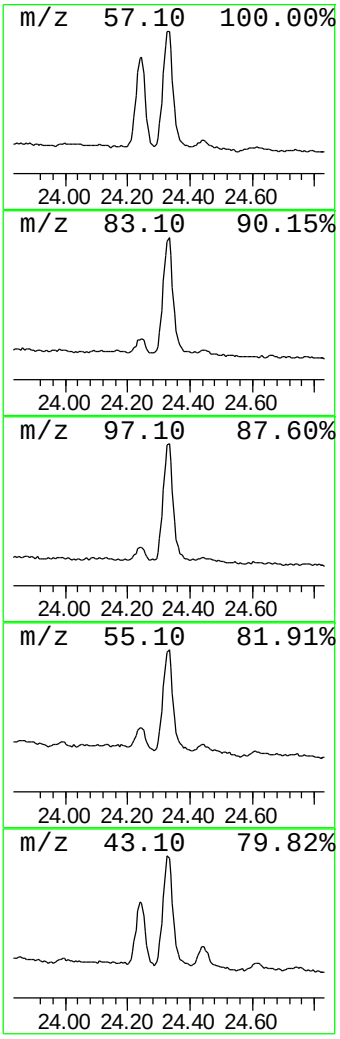
Peak Number 13 (DEL) Alkane: Cyclic-24.33 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.33	25.59 ng/ul	7755190	Perylene-d12	23.70

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9-Nonadecene	266	C19H38	031035-07-1	93
2	1-Docosene	308	C22H44	001599-67-3	91
3	17-Pentatriacontene	491	C35H70	006971-40-0	91
4	Cyclooctacosane	392	C28H56	000297-24-5	87
5	Octacosanol	410	C28H58O	000557-61-9	83





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Operator : UM/SJ
Sample : H2841-16
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : D9WT1

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM051816.M
Quant Title : SVOA CALIBRATION

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

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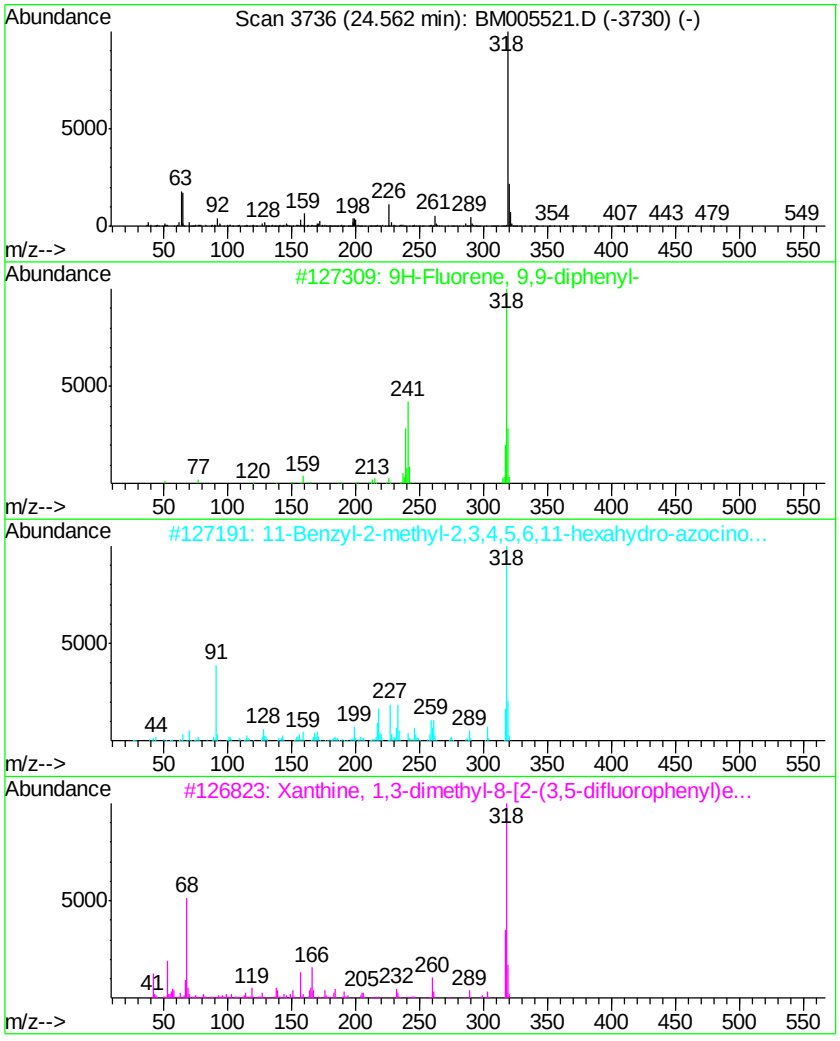
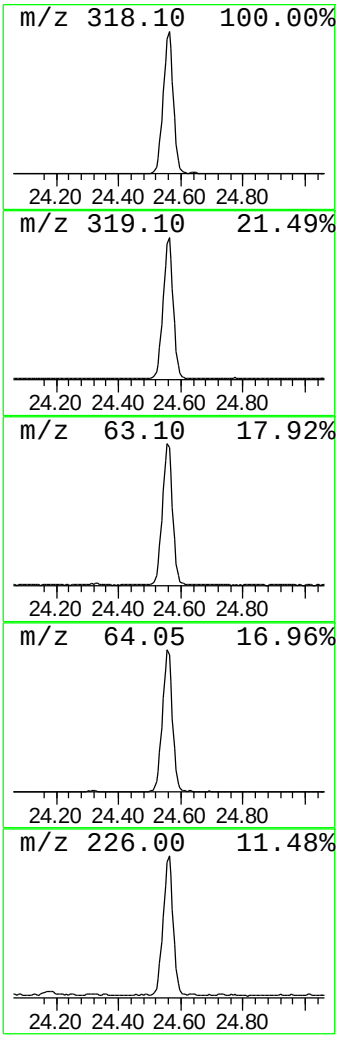
Peak Number 14 9H-Fluorene, 9,9-diphenyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.56	15.39 ng/ul	4665180	Perylene-d12	23.70

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9H-Fluorene, 9,9-diphenyl-	318	C25H18	020302-14-1	53
2	11-Benzyl-2-methyl-2,3,4,5,6,11-...	318	C21H22N2O	122531-39-9	45
3	Xanthine, 1,3-dimethyl-8-[2-(3,5...	318	C15H12F2N4O2	1000127-95-4	45
4	2,4-Diamino-5,6-diphenylthieno[2...	318	C18H14N4S	042160-23-6	42
5	Thiazole, 4-(3-fluoro-4-methoxyp...	318	C16H12F2N2OS	1000264-30-7	42





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Sample : H2841-16
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : D9WT1

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM051816.M
Quant Title : SVOA CALIBRATION

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

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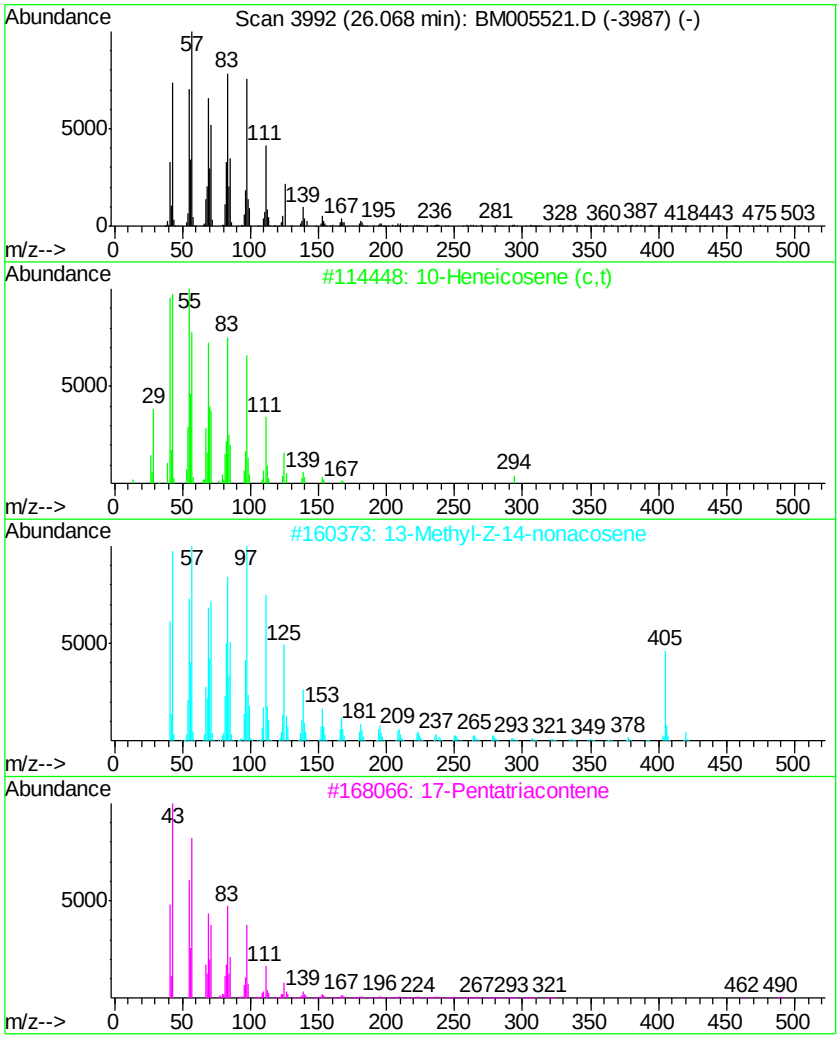
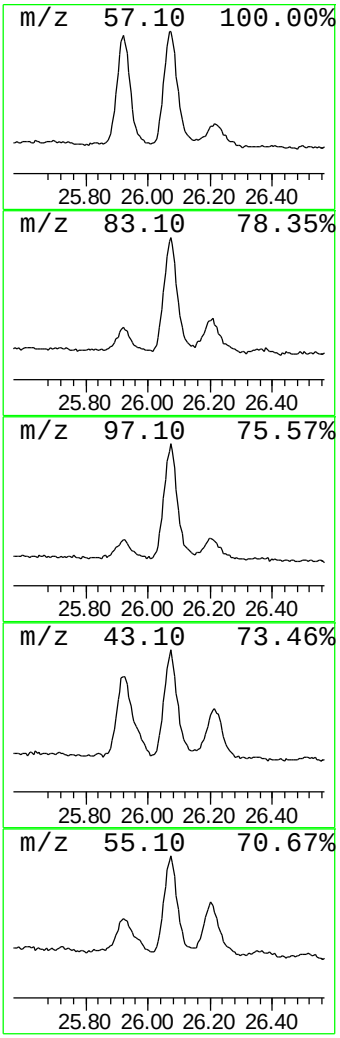
Peak Number 15 (DEL) Alkane: Cyclic-26.07 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.07	18.21 ng/ul	5520040	Perylene-d12	23.70

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	10-Heneicosene (c,t)	294	C21H42	095008-11-0	93
2	13-Methyl-Z-14-nonacosene	420	C30H60	1000131-19-0	92
3	17-Pentatriacontene	491	C35H70	006971-40-0	91
4	1-Triacontanol	438	C30H62O	000593-50-0	89
5	Pentafluoropropionic acid, octad...	416	C21H37F5O2	1000280-07-7	80





Tentatively Identified Compound (LSC) summary

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM051916\
Data File : BM005521.D
Acq On : 20 May 2016 05:11
Operator : UM/SJ
Sample : H2841-16
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : D9WT1

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM051816.M
Quant Title : SVOA CALIBRATION

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

					Internal Standard		
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
Benzenamine, 3-me...	8.75	2.0	ng/ul	100144	1	7.84	978782	20.0
Benzenamine, 2-ch...	12.22	5.0	ng/ul	414234	2	10.63	1647990	20.0
Benzenamine, 2-me...	16.62	4.4	ng/ul	817347	4	17.21	3724730	20.0
(DEL) Alkane: Cyc...	17.59	3.0	ng/ul	566144	4	17.21	3724730	20.0
Tetradecanoic acid	18.11	5.0	ng/ul	923980	4	17.21	3724730	20.0
9,10-Anthracenedione	18.63	4.3	ng/ul	801943	4	17.21	3724730	20.0
(DEL) Alkane: Cyc...	18.96	6.7	ng/ul	1241360	4	17.21	3724730	20.0
7H-Dibenzo(a,g)ca...	19.30	3.1	ng/ul	581141	4	17.21	3724730	20.0
Octadecanoic acid	19.39	3.1	ng/ul	728060	5	21.39	4638880	20.0
7H-Benz[de]anthra...	21.60	6.8	ng/ul	1570980	5	21.39	4638880	20.0
3-Thiocyanatocarb...	21.71	12.5	ng/ul	2897930	5	21.39	4638880	20.0
(DEL) Alkane: Cyc...	24.33	25.6	ng/ul	7755190	6	23.70	6061290	20.0
9H-Fluorene, 9,9-...	24.56	15.4	ng/ul	4665180	6	23.70	6061290	20.0
(DEL) Alkane: Cyc...	26.07	18.2	ng/ul	5520040	6	23.70	6061290	20.0