

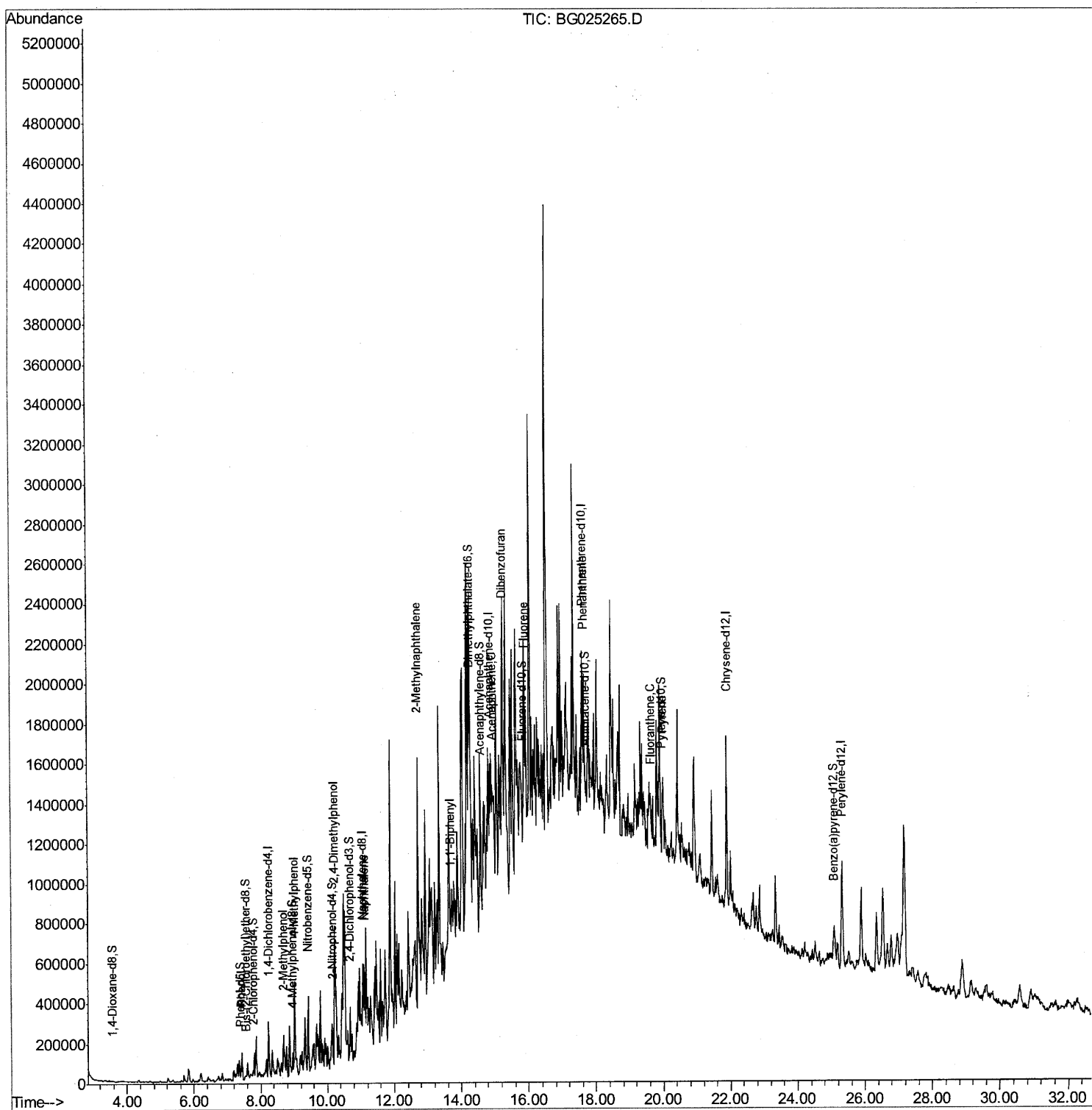
Data Path : Z:\HPCHEM1\BNA_G\DATA\BG121616\
Data File : BG025265.D
Acq On : 17 Dec 2016 8:26
Operator : UM/SJ
Sample : H5995-15DL 5X
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampled :
DA883DL

Manual Integrations
APPROVED

umangi
12/20/2016 10:49:09 AM

Quant Time: Dec 18 01:03:41 2016
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG113016.M
Quant Title : SVOA CALIBRATION
QLast Update : Sun Dec 18 00:39:32 2016
Response via : Initial Calibration



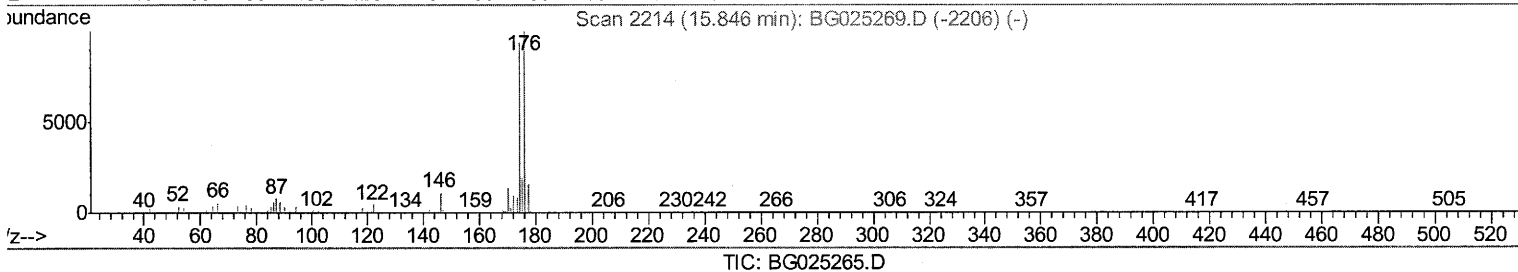
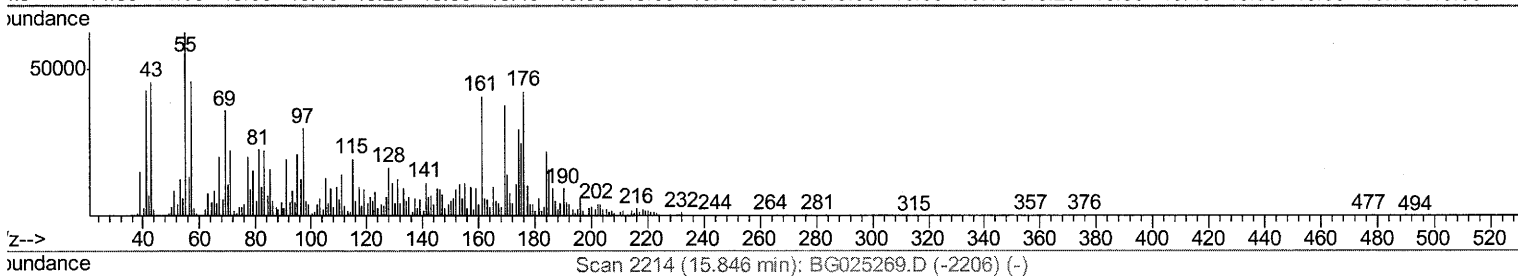
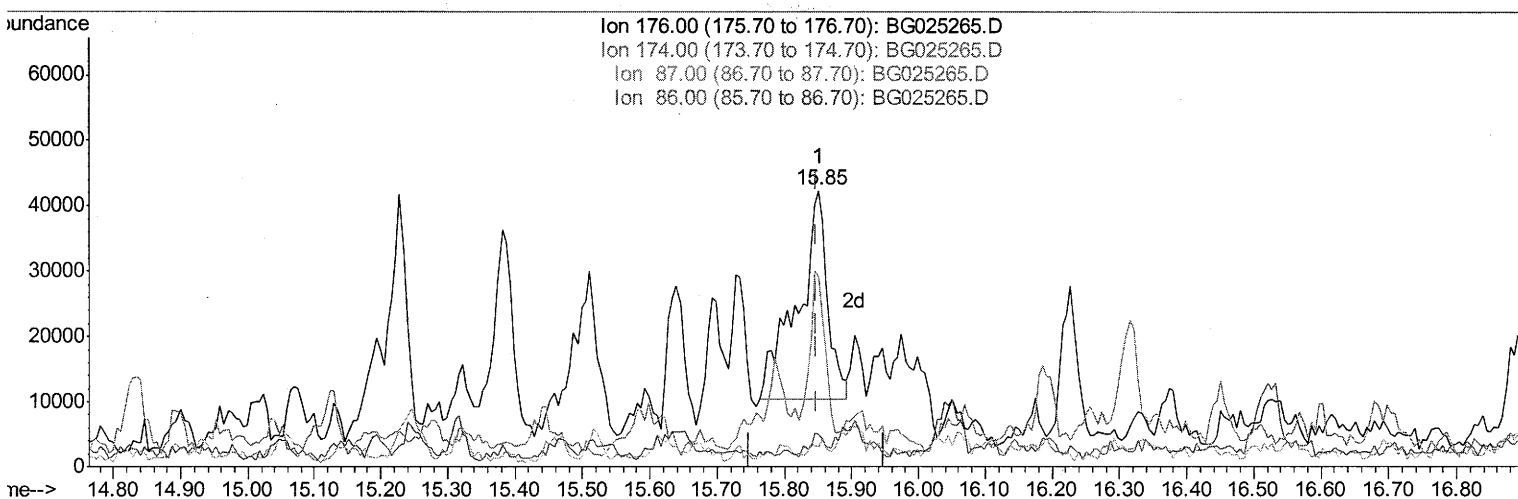
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(57) Fluorene-d10 (S)

15.851min (+0.003) 7.40ng/ul

response 100578

Ion	Exp%	Act%
176.00	100	100
174.00	92.80	69.23#
87.00	9.40	7.22#
86.00	8.20	12.04#

Quantitation Report (Qedit)

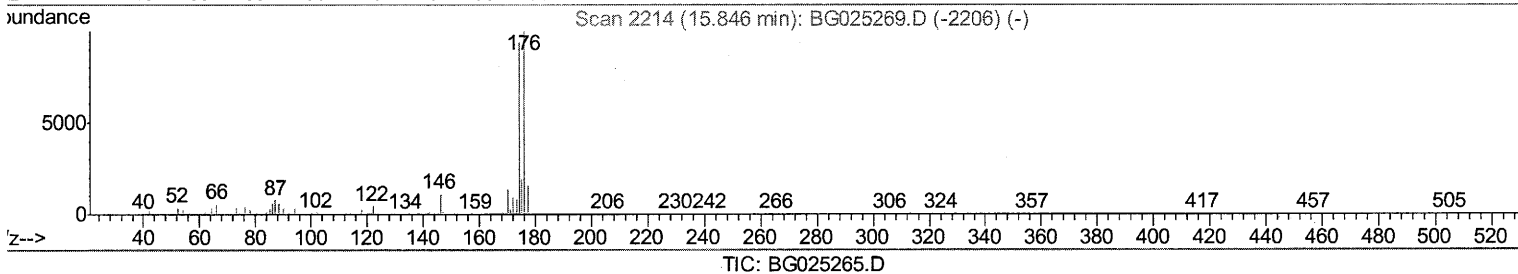
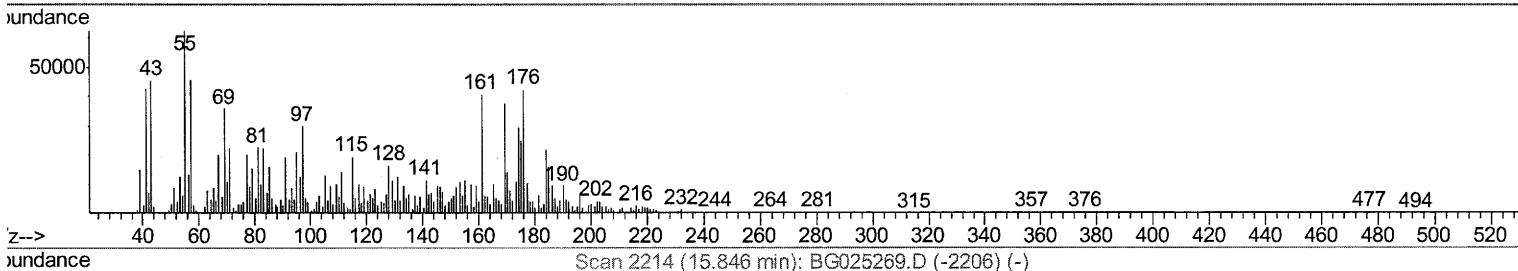
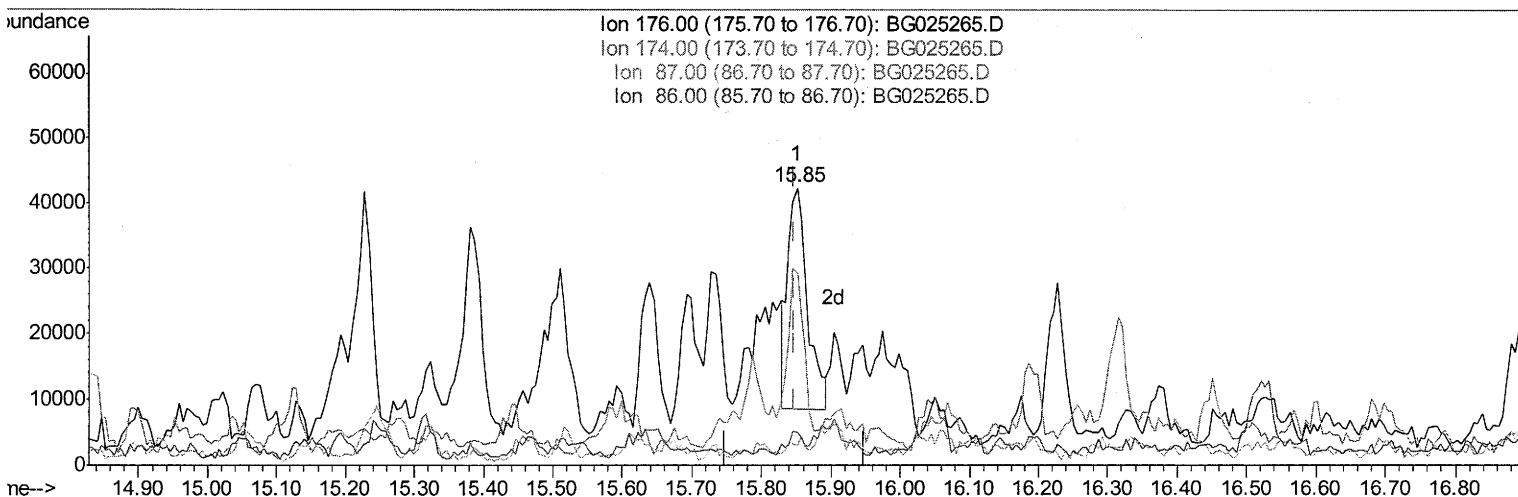
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(57) Fluorene-d10 (S)

15.851min (+0.003) 4.95ng/ul m

response 67384

Ion	Exp%	Act%
176.00	100	100
174.00	92.80	69.23#
87.00	9.40	7.22#
86.00	8.20	12.04#

Um
 12/17/16

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.23	152	71562	20.00	ng/ul	0.00
18) Naphthalene-d8	11.06	136	269305	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.86	164	210059	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.61	188	435137	20.00	ng/ul	0.00
75) Chrysene-d12	21.90	240	572226	20.00	ng/ul	0.00
83) Perylene-d12	25.33	264	600657	20.00	ng/ul	0.00

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.58	96	1751	1.26	ng/uL	0.00
5) Phenol-d5	7.40	99	17478	2.83	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.55	67	12436	3.26	ng/ul	0.00
9) 2-Chlorophenol-d4	7.77	132	12252	2.84	ng/ul	0.00
13) 4-Methylphenol-d8	8.95	113	14329	2.78	ng/ul	0.00
19) Nitrobenzene-d5	9.42	128	7947	4.17	ng/ul	0.00
22) 2-Nitrophenol-d4	10.15	143	7259	3.05	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.69	165	14627	3.20	ng/ul	0.00
29) 4-Chloroaniline-d4	0.00	131	0d	0.00	ng/ul	
43) Dimethylphthalate-d6	14.25	166	48546	3.36	ng/ul	0.00
46) Acenaphthylene-d8	14.56	160	64012	4.04	ng/ul	0.00
51) 4-Nitrophenol-d4	0.00	143	0d	0.00	ng/ul	
57) Fluorene-d10	15.85	176	67384m	4.95	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	0.00	200	0d	0.00	ng/ul	
70) Anthracene-d10	17.71	188	71833	3.91	ng/ul	0.00
76) Pyrene-d10	19.98	212	89752	3.81	ng/ul	0.00
87) Benzo(a)pyrene-d12	25.09	264	83421	3.43	ng/ul	0.00

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Target Compounds				Qvalue		
6) Phenol	7.43	94	51721	8.16	ng/ul	98
11) 2-Methylphenol	8.68	108	62257	13.34	ng/ul	85
16) 4-Methylphenol	9.02	108	164846	31.76	ng/ul	92
24) 2,4-Dimethylphenol	10.22	107	282120	51.15	ng/ul	97
28) Naphthalene	11.12	128	260434	20.80	ng/ul	99
34) 2-Methylnaphthalene	12.70	142	571138	57.81	ng/ul	94
40) 1,1'-Biphenyl	13.69	154	97579	7.66	ng/ul#	94
49) Acenaphthene	14.93	153	73058	6.93	ng/ul#	77
53) Dibenzofuran	15.26	168	126465	7.59	ng/ul	90
58) Fluorene	15.90	166	204329	15.05	ng/ul#	98
69) Phenanthrene	17.65	178	438150	21.10	ng/ul	95
74) Fluoranthene	19.64	202	66291	2.69	ng/ul#	90
77) Pyrene	20.01	202	105629	3.84	ng/ul#	90

(#) = qualifier out of range (m) = manual integration (+) = signals summed