

Quantitation Report (Qedit)

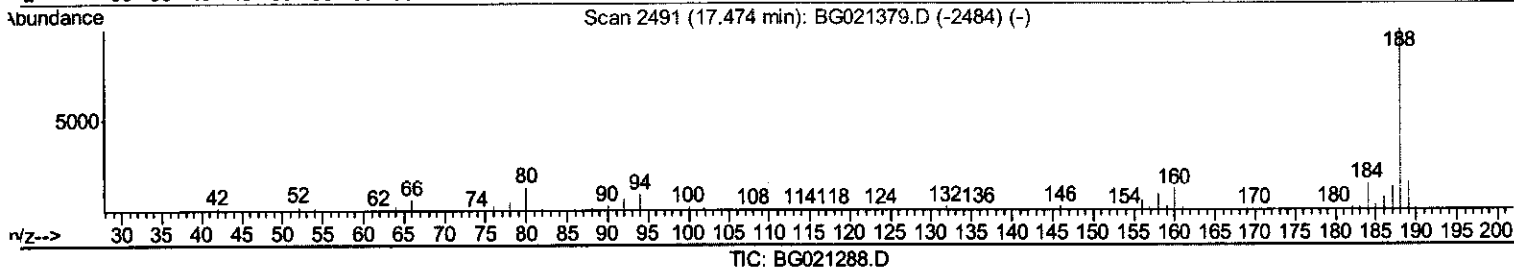
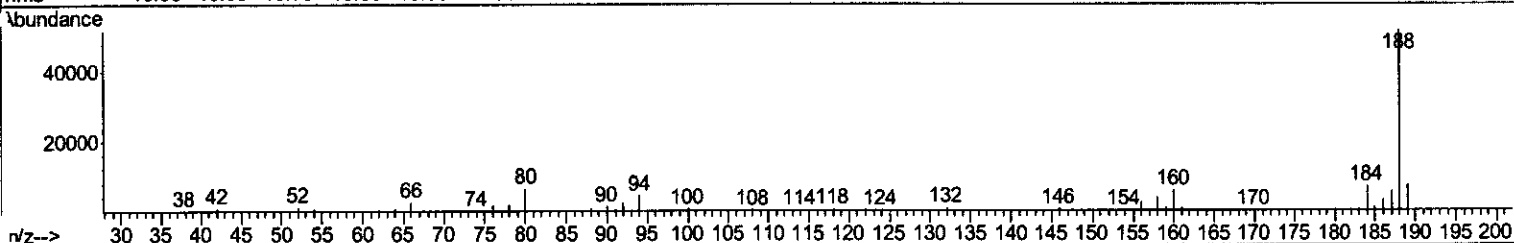
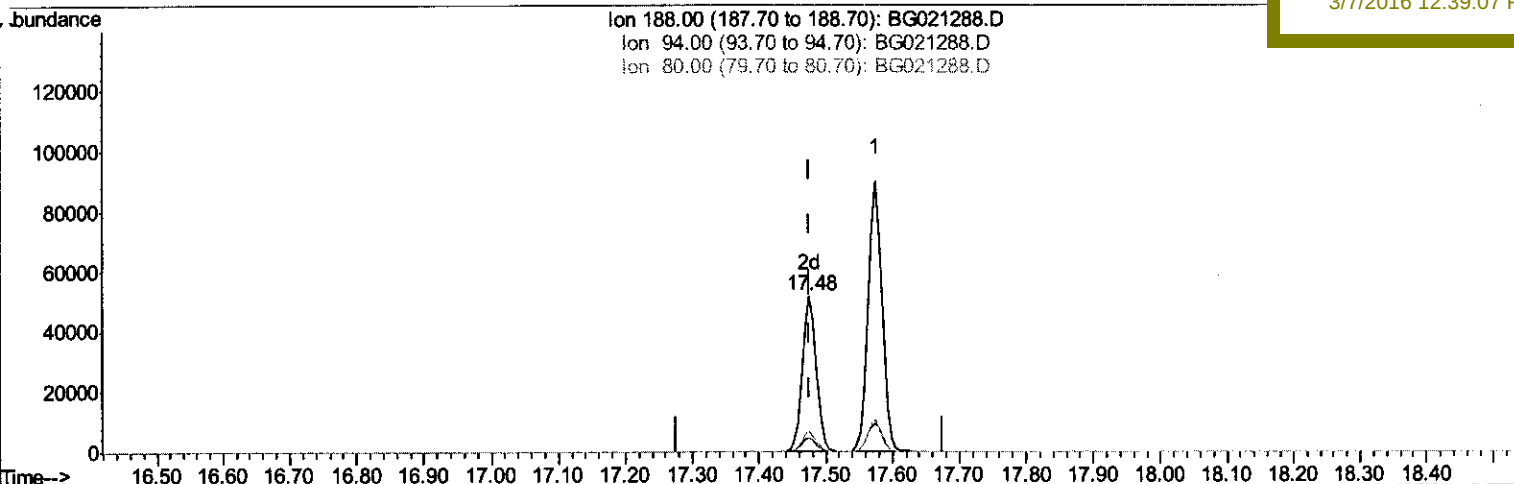
Data Path : Z:\HPCHEM1\BNA G\Data\BG030416\
 Data File : BG021288.D
 Acq On : 5 Mar 2016 2:52
 Operator : UM/SJ
 Sample : H1626-17
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Quant Time: Mar 05 06:01:26 2016
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG030316.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Mar 05 05:52:35 2016
 Response via : Initial Calibration

Instrument :
 BNA_G
 ClientSampleId :
 JHFK2

Manual Integrations
 APPROVED

Sohil
 3/7/2016 12:39:07 PM



(62) Phenanthrene-d10 (I)

17.475min (-0.001) 20.00ng/ul m

response 76052

Ion	Exp%	Act%
188.00	100	100
94.00	9.30	9.14
80.00	10.60	12.72#
0.00	0.00	0.00

19.7
 03/04/16

Quantitation Report (Qedit)

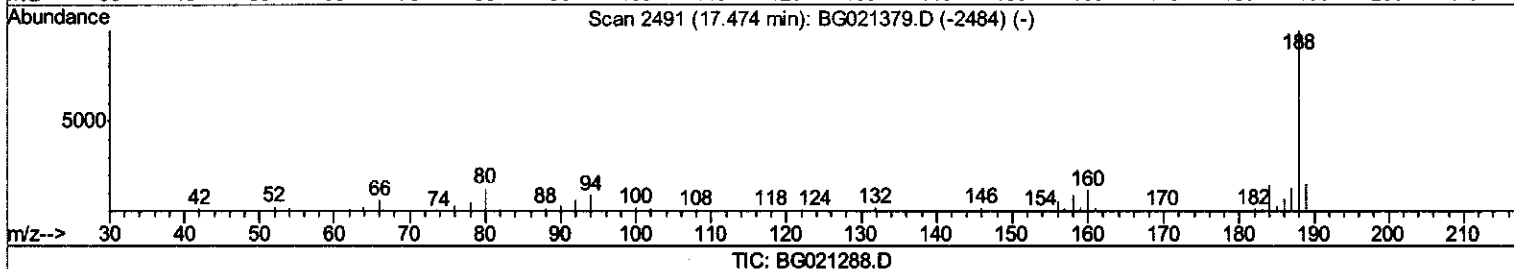
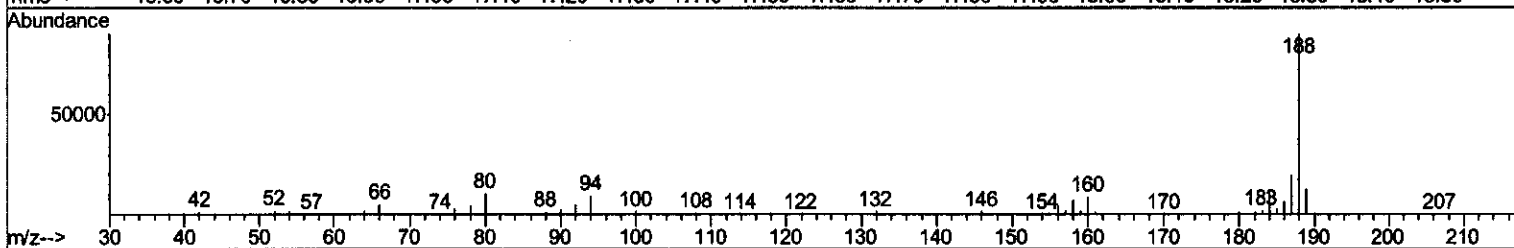
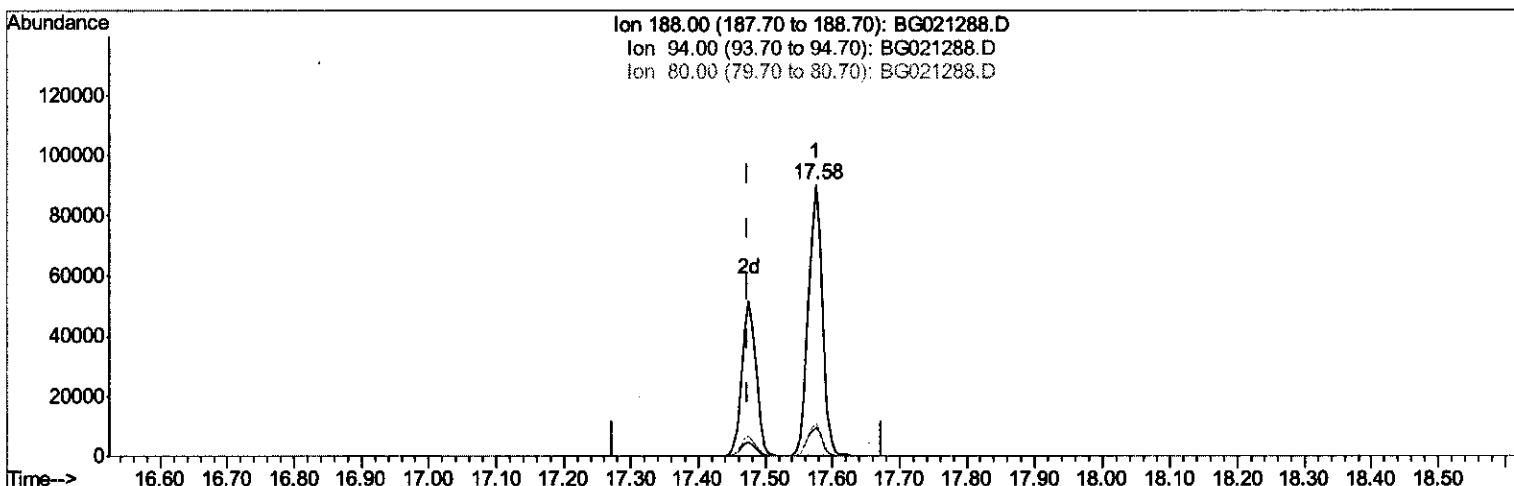
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Manual Integrations
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Sohil
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Quant Time: Mar 10 04:52:51 2016
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG030316.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Mar 09 02:03:52 2016
 Response via : Initial Calibration



(62) Phenanthrene-d10 (I)

17.575min (+0.101) 20.00ng/ul

response 132099

Ion	Exp%	Act%
188.00	100	100
94.00	9.30	10.61
80.00	10.60	11.81
0.00	0.00	0.00

Quantitation Report (LSC Reviewed)

Data Path : Z:\HPCHEM1\BNA G\Data\BG030416\
 Data File : BG021288.D
 Acq On : 5 Mar 2016 2:52
 Operator : UM/SJ
 Sample : H1626-17
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
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 ClientSampleId :
 JHFK2

Manual Integrations
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Sohil
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Quant Time: Mar 05 06:01:26 2016
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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.13	152	10514	20.00	ng/ul	0.00
18) Naphthalene-d8	10.94	136	52133	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.74	164	31649	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.48	188	76052m	20.00	ng/ul	0.00
78) Chrysene-d12	21.74	240	79713	20.00	ng/ul	0.00
86) Perylene-d12	25.00	264	74846	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.53	96	1505	5.81	ng/uL	0.00
5) Phenol-d5	7.28	99	35738	32.09	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.45	67	25198	33.93	ng/ul	0.00
9) 2-Chlorophenol-d4	7.66	132	26281	33.83	ng/ul	0.00
13) 4-Methylphenol-d8	8.83	113	29278	33.30	ng/ul	0.00
19) Nitrobenzene-d5	9.29	128	13823	32.73	ng/ul	0.00
22) 2-Nitrophenol-d4	10.01	143	14844	32.97	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.55	165	25493	31.69	ng/ul	0.00
29) 4-Chloroaniline-d4	11.07	131	39341	38.56	ng/ul	0.00
44) Dimethylphthalate-d6	14.14	166	92457	34.86	ng/ul	0.00
47) Acenaphthylene-d8	14.44	160	114194	34.60	ng/ul	0.00
52) 4-Nitrophenol-d4	14.92	143	17637	34.12	ng/ul	0.00
58) Fluorene-d10	15.72	176	82527	35.43	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.84	200	15781	28.91	ng/ul	0.00
71) Anthracene-d10	17.58	188	132099	34.93	ng/ul	0.00
79) Pyrene-d10	19.85	212	150056	37.19	ng/ul	0.00
90) Benzo(a)pyrene-d12	24.77	264	142898	35.37	ng/ul	0.00

Target Compounds

					Ovalue
45) Dimethylphthalate	14.19	163	51078	17.90 ng/ul#	98

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

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