

Quantitation Report (QT/LSC Reviewed)

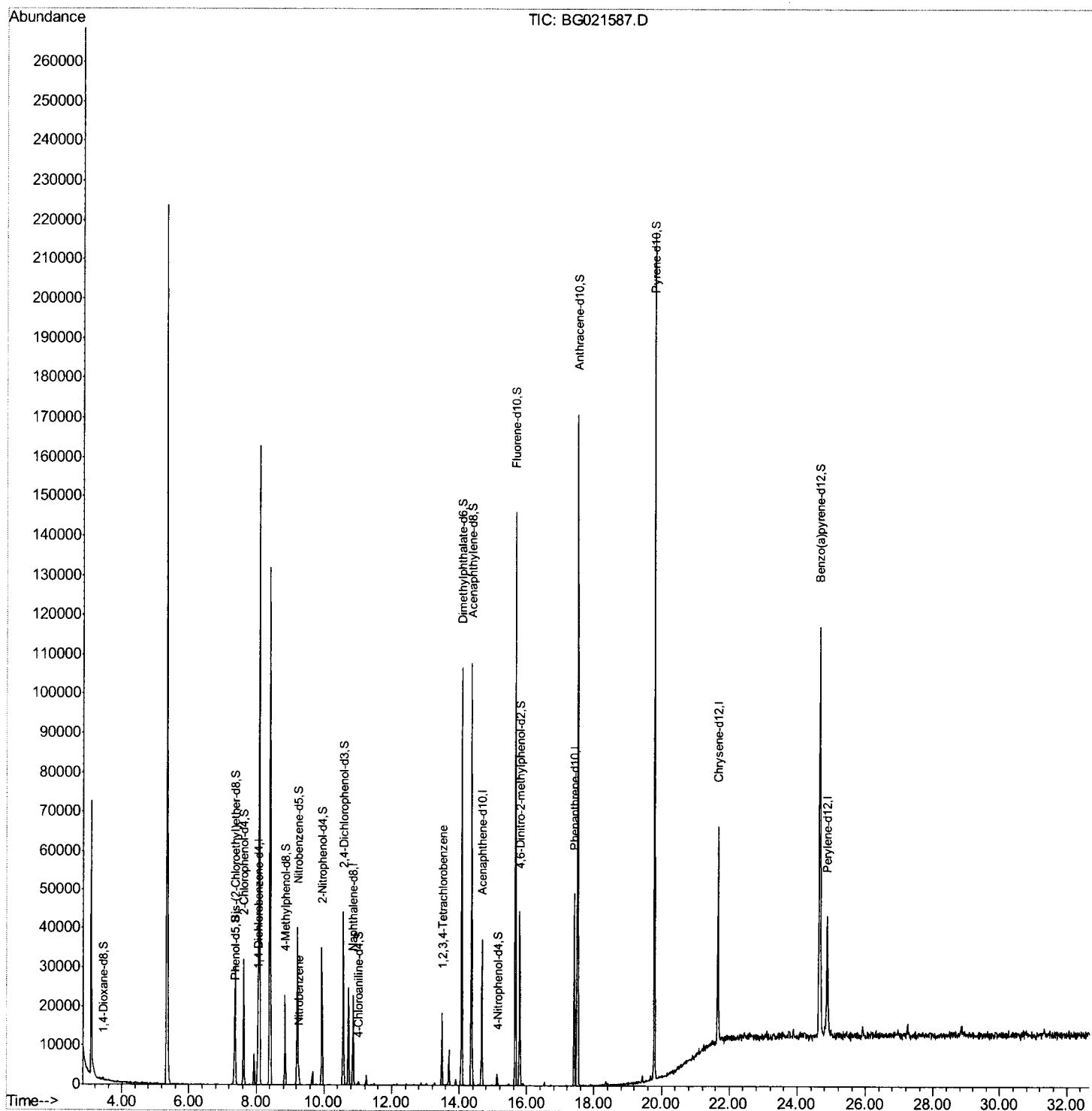
Data Path : Z:\HPCHEM1\BNA_G\DATA\BG033016\
 Data File : BG021587.D
 Acq On : 30 Mar 2016 19:04
 Operator : UM/SJ
 Sample : H1935-13
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 C0AJ4

Manual Integrations
 APPROVED

Sohil
 3/31/2016 1:23:59 PM

Quant Time: Mar 31 11:20:28 2016
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG031616.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Mar 31 02:07:33 2016
 Response via : Initial Calibration



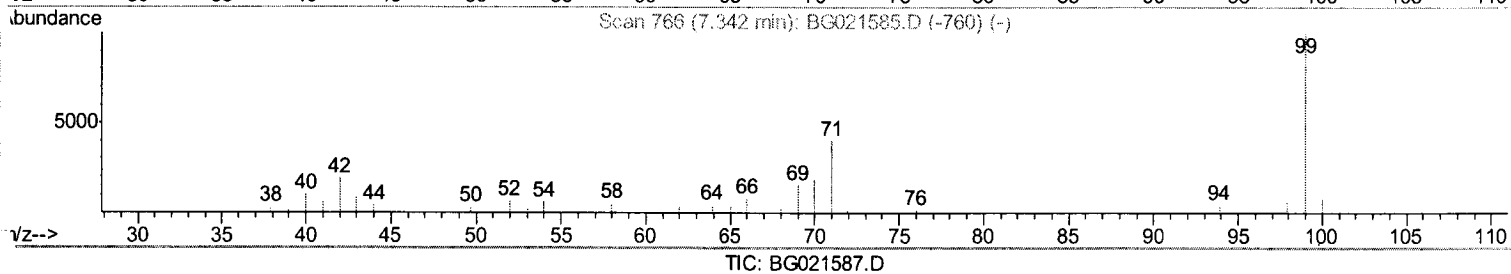
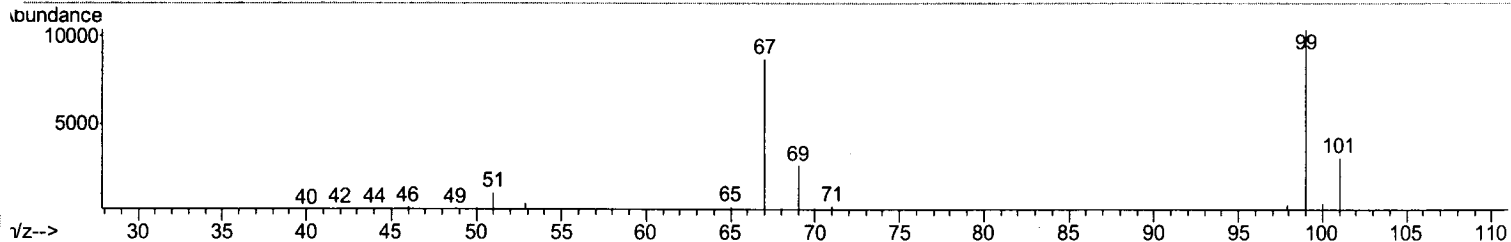
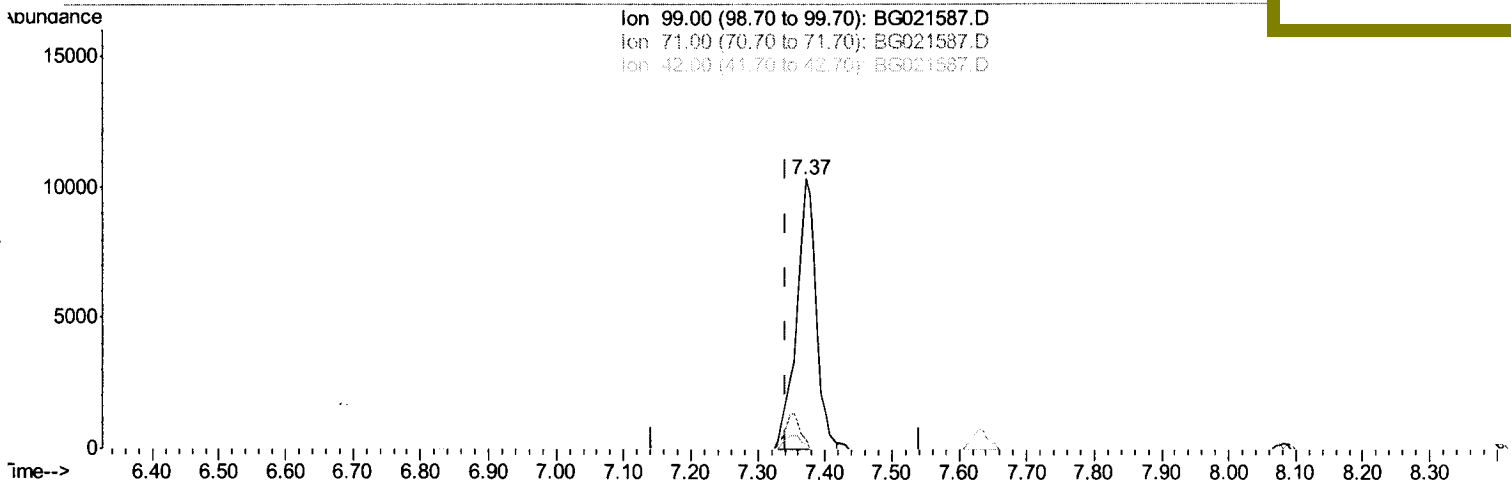
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TIC: BG021587.D

(5) Phenol-d5 (S)

7.371min (+0.030) 46.21ng/ul

response 20868

Ion	Exp%	Act%
99.00	100	100
71.00	46.50	3.52#
42.00	19.40	1.72#
0.00	0.00	0.00

Quantitation Report (Qedit)

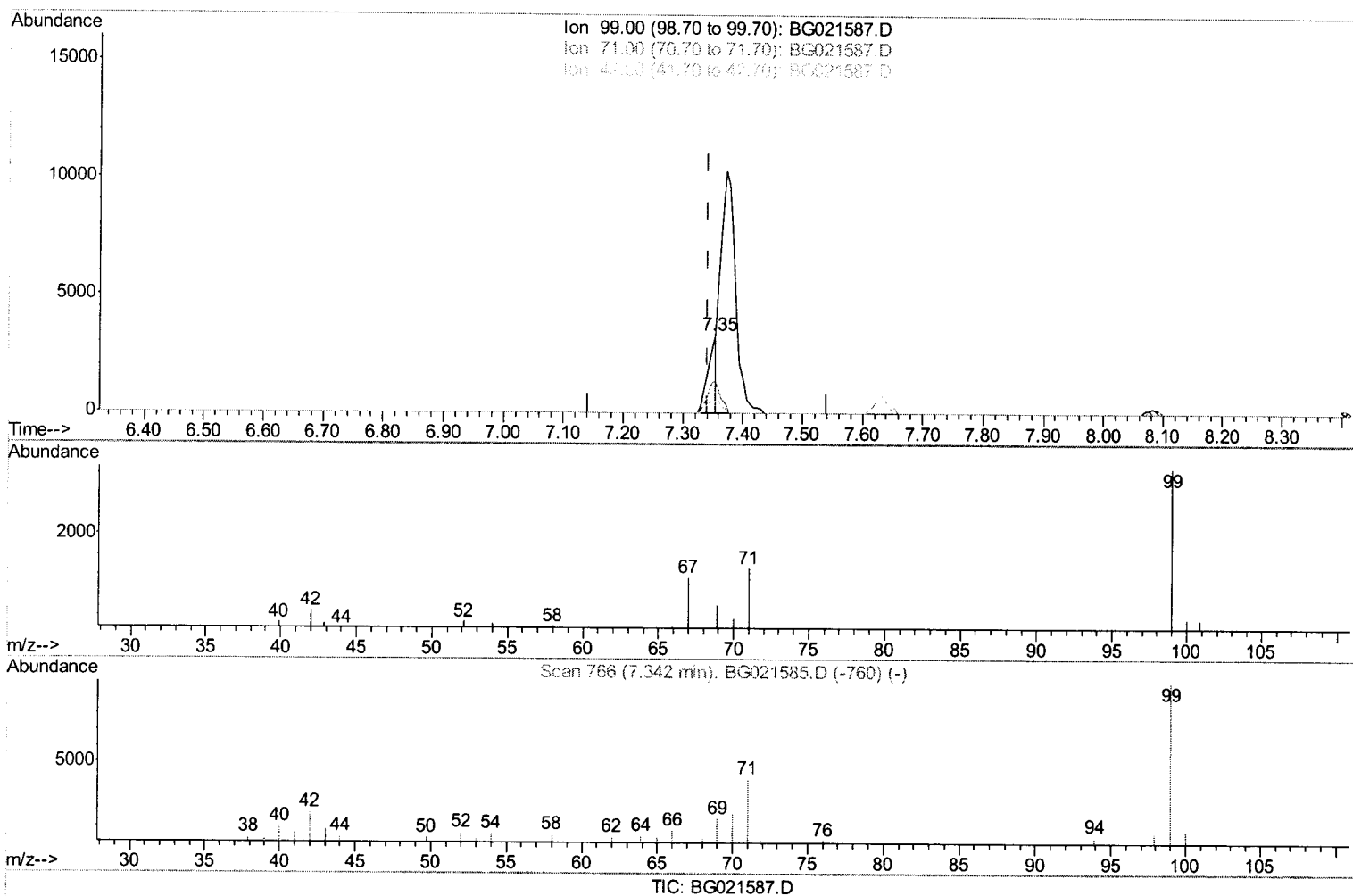
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Manual Integrations
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(5) Phenol-d5 (S)

7.354min (+0.012) 7.12ng/ul m

response 3214

> U.M
 04/06/16

Ion	Exp%	Act%
99.00	100	100
71.00	46.50	40.47
42.00	19.40	14.88#
0.00	0.00	0.00

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Manual Integrations
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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.05	152	4777	20.00	ng/ul	0.00
18) Naphthalene-d8	10.86	136	19988	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.67	164	12433	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.41	188	31450	20.00	ng/ul	0.00
78) Chrysene-d12	21.67	240	35042	20.00	ng/ul	0.00
86) Perylene-d12	24.87	264	33186	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.45	96	140	1.31	ng/uL	0.00
5) Phenol-d5	7.35	99	3214m	7.12	ng/ul	0.01
7) Bis-(2-Chloroethyl)ether-d	7.37	67	14550	54.38	ng/ul	0.00
9) 2-Chlorophenol-d4	7.63	132	14821	43.21	ng/ul	0.00
13) 4-Methylphenol-d8	8.85	113	9307	24.06	ng/ul	0.00
19) Nitrobenzene-d5	9.22	128	9696	64.66	ng/ul	0.00
22) 2-Nitrophenol-d4	9.94	143	10731	62.60	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.57	165	16814	54.87	ng/ul	0.00
29) 4-Chloroaniline-d4	11.01	131	1204	3.01	ng/ul	0.00
44) Dimethylphthalate-d6	14.07	166	65057	62.84	ng/ul	0.00
47) Acenaphthylene-d8	14.36	160	80483	63.07	ng/ul	0.00
52) 4-Nitrophenol-d4	15.11	143	980	5.92	ng/ul	0.02
58) Fluorene-d10	15.66	176	58607	62.79	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.79	200	10505	50.56	ng/ul	0.00
71) Anthracene-d10	17.51	188	97618	63.34	ng/ul	0.00
79) Pyrene-d10	19.79	212	115139	70.46	ng/ul	0.00
90) Benzo(a)pyrene-d12	24.66	264	114271	74.61	ng/ul	0.00

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 04/06/16

Target Compounds

						Qvalue
20) Nitrobenzene	9.26	77	2742	6.47	ng/ul	92
73) 1,2,3,4-Tetrachlorobenzene	13.48	216	885	2.01	ng/uL	84

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