

Quantitation Report (QT Reviewed)

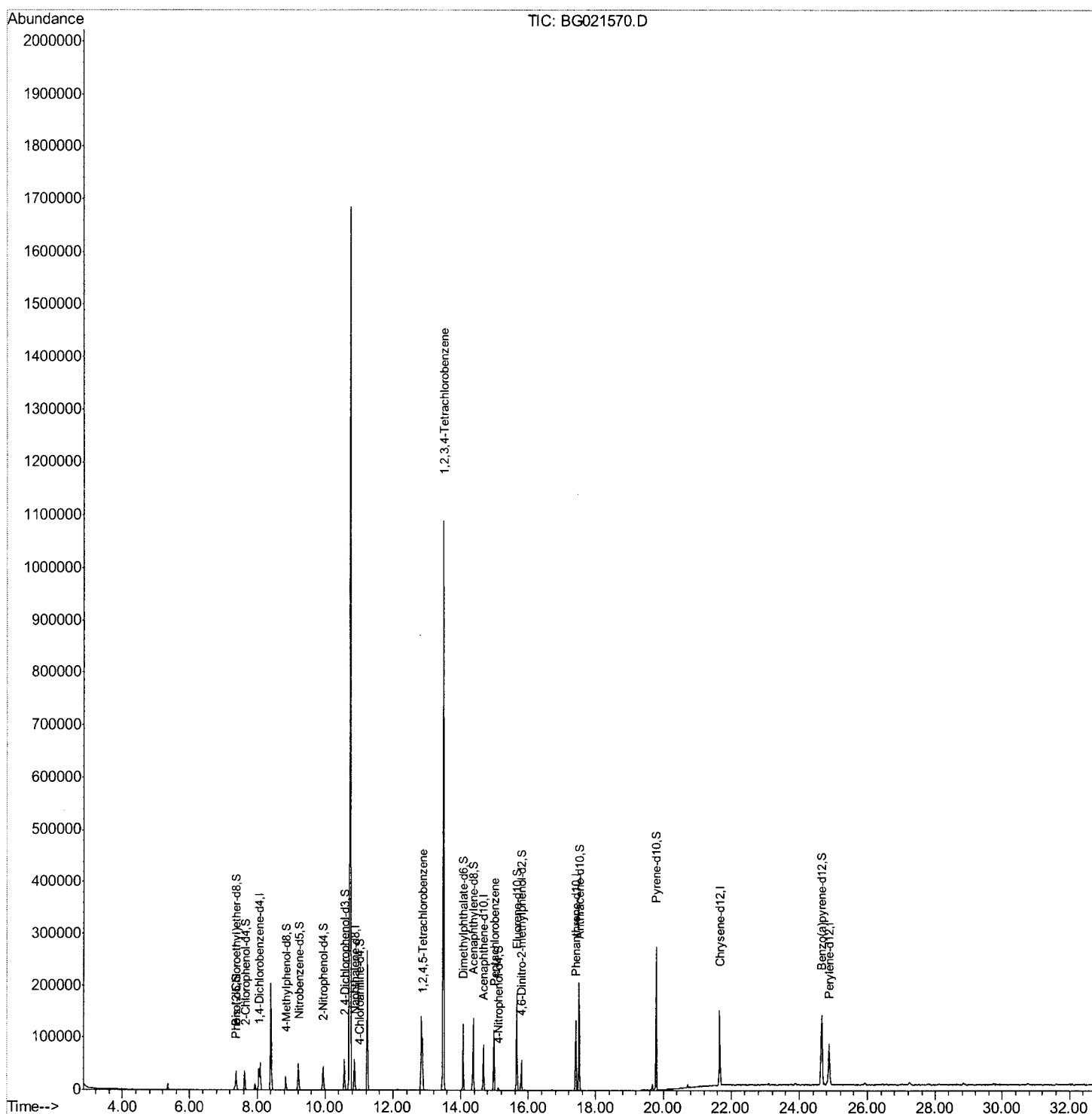
Data Path : Z:\HPCHEM1\BNA_G\DATA\BG032916\
 Data File : BG021570.D
 Acq On : 30 Mar 2016 2:51
 Operator : UM/SJ
 Sample : H1910-12
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 C0AB3

Manual Integrations
 APPROVED

Sohil
 3/30/2016 7:15:50 PM

Quant Time: Mar 30 04:55:13 2016
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG031616.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Mar 30 01:34:35 2016
 Response via : Initial Calibration



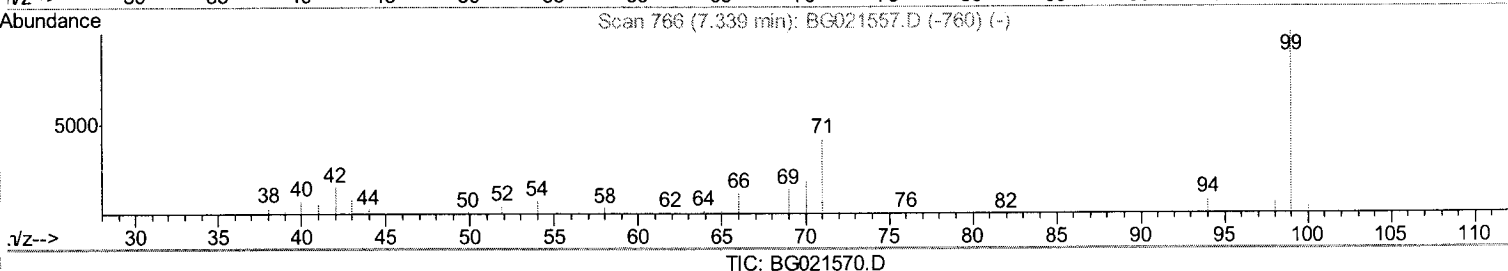
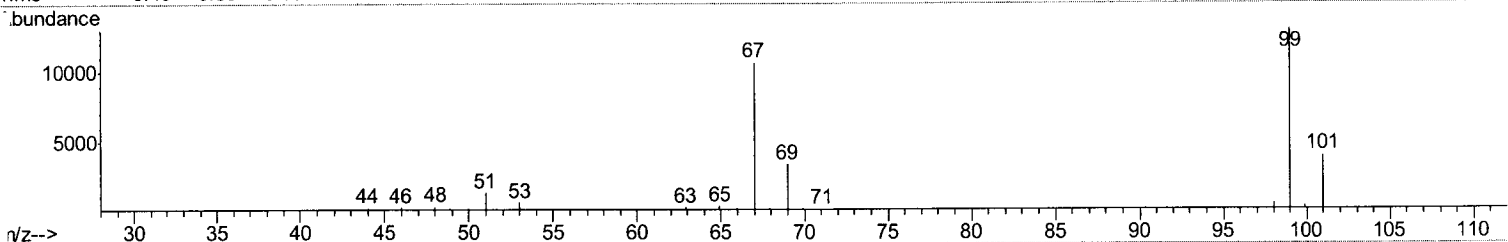
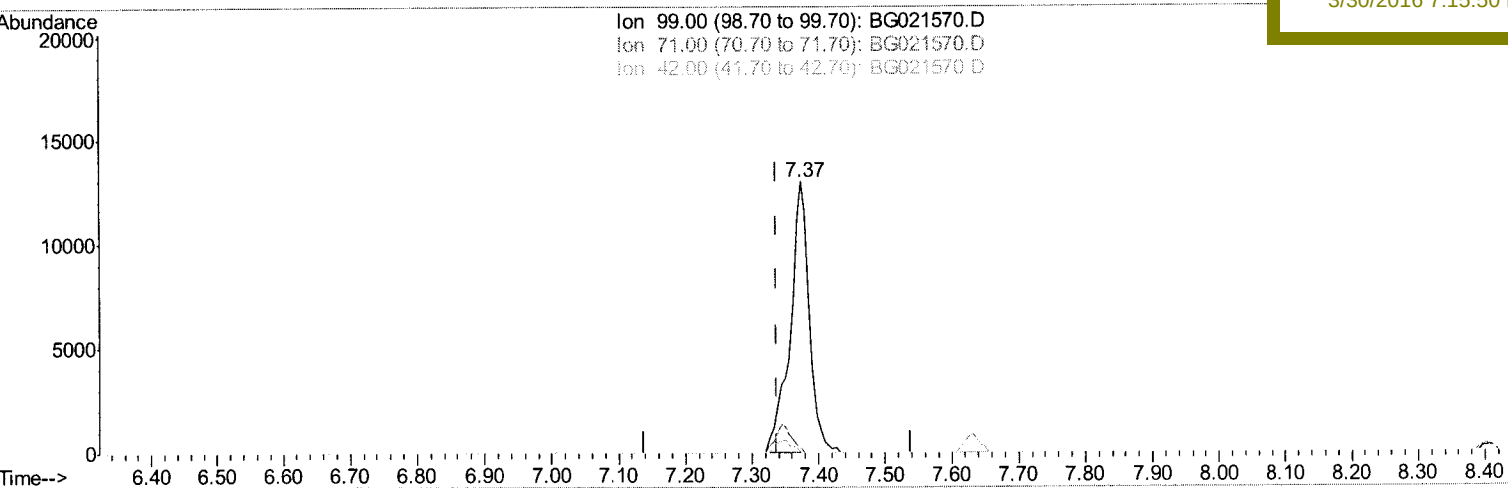
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(5) Phenol-d5 (S)

7.374min (+0.035) 25.30ng/ul

response 26112

Ion	Exp%	Act%
99.00	100	100
71.00	46.50	1.25#
42.00	19.40	0.00#
0.00	0.00	0.00

Quantitation Report (Qedit)

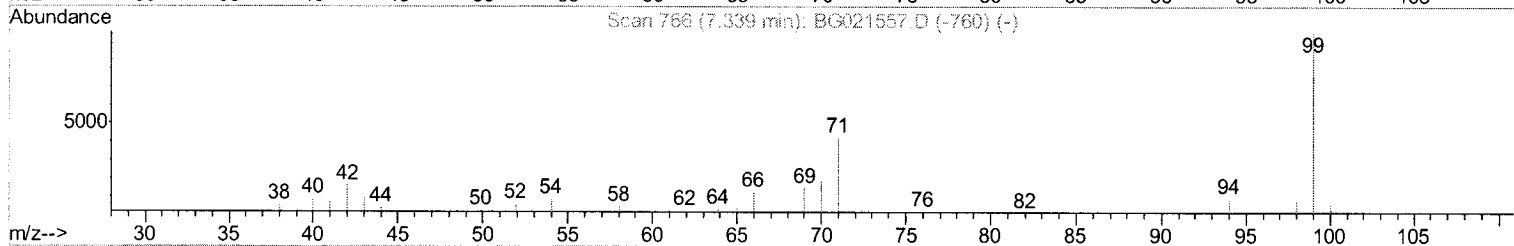
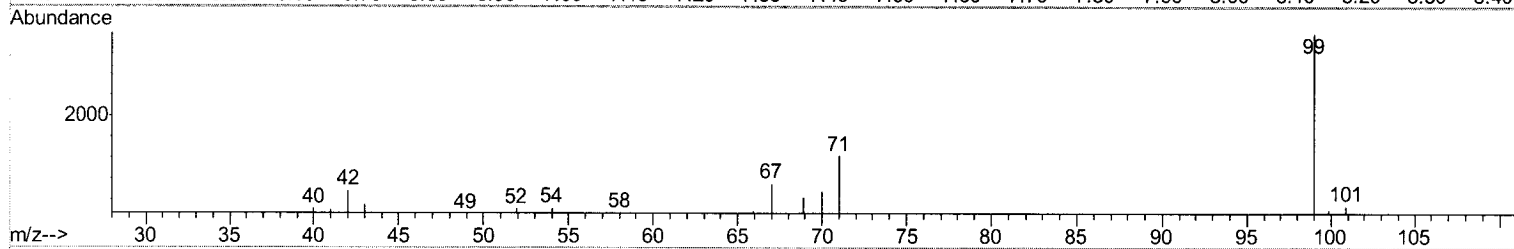
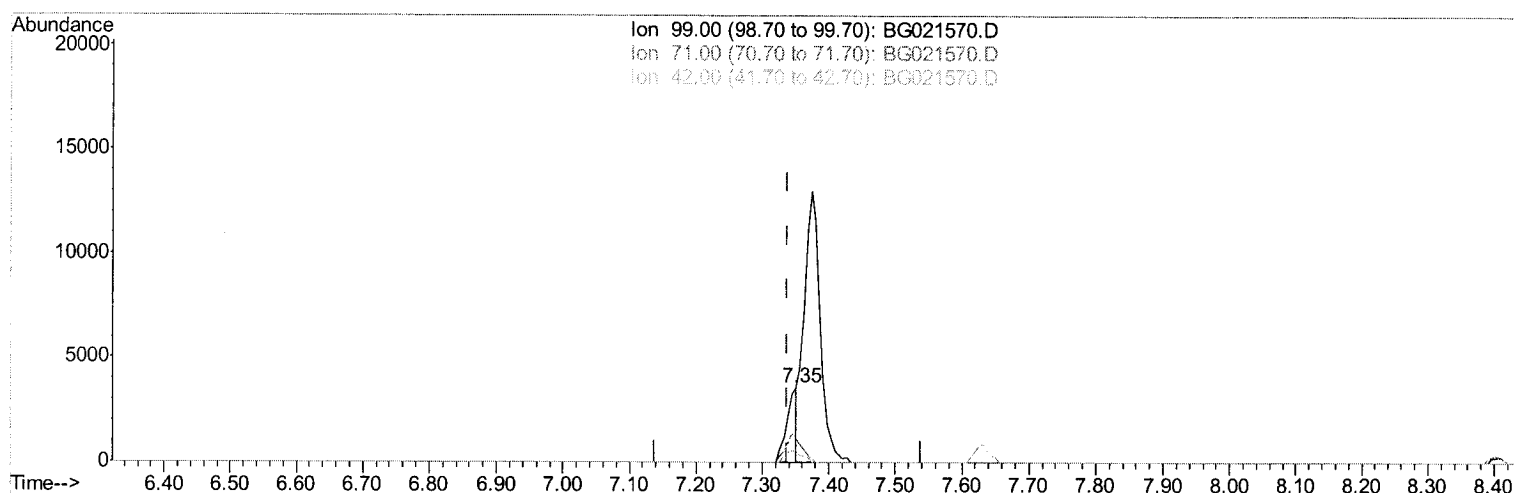
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Manual Integrations
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TIC: BG021570.D

(5) Phenol-d5 (S)

7.350min (+0.011) 3.75ng/ul m

response 3873

U.M
 04/11/16

Ion	Exp%	Act%
99.00	100	100
71.00	46.50	34.79#
42.00	19.40	16.29
0.00	0.00	0.00

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Manual Integrations
 APPROVED

Sohil
 3/30/2016 7:15:50 PM

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev (Min)
1) 1,4-Dichlorobenzene-d4	8.05	152	10916	20.00	ng/ul	0.00
18) Naphthalene-d8	10.86	136	46753	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.67	164	29628	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.41	188	76190	20.00	ng/ul	0.00
78) Chrysene-d12	21.66	240	89256	20.00	ng/ul	0.00
86) Perylene-d12	24.87	264	87640	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.44	96	176	0.72	ng/uL	0.00
5) Phenol-d5	7.35	99	3873m	3.75	ng/ul	0.01
7) Bis-(2-Chloroethyl)ether-d	7.37	67	17644	28.86	ng/ul	0.00
9) 2-Chlorophenol-d4	7.63	132	17874	22.81	ng/ul	0.00
13) 4-Methylphenol-d8	8.84	113	10989	12.43	ng/ul	0.00
19) Nitrobenzene-d5	9.22	128	12273	34.99	ng/ul	0.00
22) 2-Nitrophenol-d4	9.94	143	12997	32.42	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.56	165	20955	29.24	ng/ul	0.00
29) 4-Chloroaniline-d4	11.01	131	1206	1.29	ng/ul	0.00
44) Dimethylphthalate-d6	14.07	166	78631	31.87	ng/ul	0.00
47) Acenaphthylene-d8	14.37	160	95963	31.56	ng/ul	0.00
52) 4-Nitrophenol-d4	15.11	143	1516	3.84	ng/ul	0.01
58) Fluorene-d10	15.66	176	72204	32.46	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.79	200	13633	27.08	ng/ul	0.00
71) Anthracene-d10	17.51	188	120788	32.35	ng/ul	0.00
79) Pyrene-d10	19.79	212	145821	35.03	ng/ul	0.00
90) Benzo(a)pyrene-d12	24.65	264	147152	36.38	ng/ul	0.00

U.M
 30411116

Target Compounds						Qvalue
37) 1,2,4,5-Tetrachlorobenzene	12.87	216	24562	23.68	ng/ul	94
73) 1,2,3,4-Tetrachlorobenzene	13.48	216	344720	323.13	ng/uL	92
74) Pentachlorobenzene	14.98	250	28960	26.15	ng/uL	98

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