

Data Path : Z:\HPCHEM1\BNA G\DATA\BG032416\
 Data File : BG021482.D
 Acq On : 24 Mar 2016 12:15
 Operator : UM/SJ
 Sample : PB89111BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK11

Quant Time: Mar 25 01:03:52 2016
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG031616.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Mar 25 01:02:38 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.06	152	8402	20.00	ng/ul	0.00
18) Naphthalene-d8	10.87	136	39236	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.68	164	26801	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.42	188	67022	20.00	ng/ul	0.00
78) Chrysene-d12	21.67	240	80855	20.00	ng/ul	0.00
86) Perylene-d12	24.88	264	78236	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.46	96	1052	5.62	ng/uL	0.00
5) Phenol-d5	7.33	99	22719	28.60	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.39	67	14994	31.86	ng/ul	0.00
9) 2-Chlorophenol-d4	7.63	132	17526	29.05	ng/ul	0.00
13) 4-Methylphenol-d8	8.84	113	19551	28.74	ng/ul	0.00
19) Nitrobenzene-d5	9.23	128	9857	33.48	ng/ul	0.00
22) 2-Nitrophenol-d4	9.95	143	10908	32.42	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.56	165	18807	31.27	ng/ul	0.00
29) 4-Chloroaniline-d4	11.01	131	29757	37.84	ng/ul	0.00
44) Dimethylphthalate-d6	14.08	166	85337	38.24	ng/ul	0.00
47) Acenaphthylene-d8	14.37	160	89339	32.48	ng/ul	0.00
52) 4-Nitrophenol-d4	15.08	143	11118	31.17	ng/ul	0.00
58) Fluorene-d10	15.66	176	68489	34.04	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.79	200	11956	27.00	ng/ul	0.00
71) Anthracene-d10	17.52	188	119476	36.38	ng/ul	0.00
79) Pyrene-d10	19.79	212	140427	37.24	ng/ul	0.00
90) Benzo(a)pyrene-d12	24.66	264	146387	40.54	ng/ul	0.00

Target Compounds Ovalue

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

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