

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG111616\
 Data File : BG024798.D
 Acq On : 16 Nov 2016 11:15
 Operator : UM/SJ
 Sample : PB94740BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK40

Quant Time: Nov 17 01:28:55 2016
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG111416.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Nov 17 01:26:41 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.27	152	83937	20.00	ng/ul	0.00
18) Naphthalene-d8	11.09	136	343627	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.88	164	296698	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.62	188	706911	20.00	ng/ul	0.00
75) Chrysene-d12	21.91	240	997619	20.00	ng/ul	0.00
83) Perylene-d12	25.35	264	1033926	20.00	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	3.62	96	14711	9.08	ng/uL	0.00
5) Phenol-d5	7.40	99	279678	38.89	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.57	67	182172	40.27	ng/ul	0.00
9) 2-Chlorophenol-d4	7.78	132	203450	39.45	ng/ul	0.00
13) 4-Methylphenol-d8	8.95	113	228283	37.81	ng/ul	0.00
19) Nitrobenzene-d5	9.43	128	102806	39.51	ng/ul	0.00
22) 2-Nitrophenol-d4	10.16	143	125091	40.28	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.69	165	226231	36.59	ng/ul	0.00
29) 4-Chloroaniline-d4	11.22	131	287630	41.65	ng/ul	0.00
43) Dimethylphthalate-d6	14.27	166	874027	40.40	ng/ul	0.00
46) Acenaphthylene-d8	14.58	160	912208	38.70	ng/ul	0.00
51) 4-Nitrophenol-d4	15.06	143	141130	36.66	ng/ul	0.00
57) Fluorene-d10	15.86	176	777358	38.18	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.97	200	172214	35.44	ng/ul	0.00
70) Anthracene-d10	17.72	188	1293220	41.42	ng/ul	0.00
76) Pyrene-d10	19.99	212	1699879	39.30	ng/ul	0.00
87) Benzo(a)pyrene-d12	25.11	264	1843365	40.73	ng/ul	0.00

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

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