

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG030316\
 Data File : BG021247.D
 Acq On : 3 Mar 2016 23:35
 Operator : UM/SJ
 Sample : PB88600BL
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK00

Quant Time: Mar 04 02:40:38 2016
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG030316.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Mar 04 02:35:52 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.13	152	9714	20.00	ng/ul	0.00
18) Naphthalene-d8	10.94	136	45917	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.74	164	27864	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.48	188	66744	20.00	ng/ul	0.00
78) Chrysene-d12	21.74	240	72103	20.00	ng/ul	0.00
86) Perylene-d12	24.99	264	67052	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.53	96	1849	7.73	ng/uL	0.00
5) Phenol-d5	7.28	99	33856	32.91	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.45	67	24898	36.29	ng/ul	0.00
9) 2-Chlorophenol-d4	7.66	132	25507	35.53	ng/ul	0.00
13) 4-Methylphenol-d8	8.82	113	28297	34.83	ng/ul	0.00
19) Nitrobenzene-d5	9.30	128	13846	37.22	ng/ul	0.00
22) 2-Nitrophenol-d4	10.02	143	14912	37.61	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.55	165	23865	33.68	ng/ul	0.00
29) 4-Chloroaniline-d4	11.07	131	38701	43.07	ng/ul	0.00
44) Dimethylphthalate-d6	14.14	166	90270	38.66	ng/ul	0.00
47) Acenaphthylene-d8	14.43	160	107522	37.00	ng/ul	0.00
52) 4-Nitrophenol-d4	14.92	143	15979	35.11	ng/ul	0.00
58) Fluorene-d10	15.73	176	77450	37.77	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.84	200	15334	32.01	ng/ul	0.00
71) Anthracene-d10	17.58	188	124502	37.51	ng/ul	0.00
79) Pyrene-d10	19.84	212	142143	38.94	ng/ul	0.00
90) Benzo(a)pyrene-d12	24.77	264	134975	37.29	ng/ul	0.00

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

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