

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG032116\
 Data File : BG021448.D
 Acq On : 21 Mar 2016 17:47
 Operator : UM/SJ
 Sample : PB89035BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK35

Quant Time: Mar 22 01:44:11 2016
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG031616.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Mar 22 01:42:49 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.07	152	9557	20.00	ng/ul	0.00
18) Naphthalene-d8	10.88	136	45961	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.69	164	31544	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.43	188	82072	20.00	ng/ul	0.00
78) Chrysene-d12	21.68	240	93558	20.00	ng/ul	0.00
86) Perylene-d12	24.91	264	91819	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.47	96	1275	5.99	ng/uL	0.00
5) Phenol-d5	7.33	99	25643	28.38	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.40	67	16576	30.97	ng/ul	0.00
9) 2-Chlorophenol-d4	7.64	132	21232	30.94	ng/ul	0.00
13) 4-Methylphenol-d8	8.85	113	23523	30.40	ng/ul	0.00
19) Nitrobenzene-d5	9.25	128	11823	34.29	ng/ul	0.00
22) 2-Nitrophenol-d4	9.96	143	13248	33.61	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.57	165	23221	32.96	ng/ul	0.00
29) 4-Chloroaniline-d4	11.02	131	33585	36.46	ng/ul	0.00
44) Dimethylphthalate-d6	14.09	166	95923	36.52	ng/ul	0.00
47) Acenaphthylene-d8	14.38	160	109396	33.79	ng/ul	0.00
52) 4-Nitrophenol-d4	15.07	143	13104	31.21	ng/ul	0.00
58) Fluorene-d10	15.68	176	80719	34.09	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.80	200	10692	19.72	ng/ul	0.00
71) Anthracene-d10	17.53	188	135212	33.62	ng/ul	0.00
79) Pyrene-d10	19.80	212	152871	35.04	ng/ul	0.00
90) Benzo(a)pyrene-d12	24.68	264	160928	37.97	ng/ul	0.00

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

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