

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG030316\
 Data File : BG021248.D
 Acq On : 4 Mar 2016 00:14
 Operator : UM/SJ
 Sample : H1626-08
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 JHFJ5

Quant Time: Mar 04 02:40:45 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	10205	20.00	ng/ul	0.00
18) Naphthalene-d8	10.94	136	49175	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.74	164	29060	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.47	188	65817	20.00	ng/ul	0.00
78) Chrysene-d12	21.74	240	72182	20.00	ng/ul	0.00
86) Perylene-d12	24.99	264	68242	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.54	96	287	1.14	ng/uL	0.00
5) Phenol-d5	7.28	99	8068	7.46	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.45	67	24666	34.22	ng/ul	0.00
9) 2-Chlorophenol-d4	7.66	132	21001	27.85	ng/ul	0.00
13) 4-Methylphenol-d8	8.83	113	15380	18.02	ng/ul	0.00
19) Nitrobenzene-d5	9.30	128	13887	34.86	ng/ul	0.00
22) 2-Nitrophenol-d4	10.02	143	15069	35.48	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.55	165	22466	29.61	ng/ul	0.00
29) 4-Chloroaniline-d4	0.00	131	0	0.00	ng/ul	
44) Dimethylphthalate-d6	14.14	166	87396	35.89	ng/ul	0.00
47) Acenaphthylene-d8	14.44	160	110892	36.59	ng/ul	0.00
52) 4-Nitrophenol-d4	14.92	143	3276	6.90	ng/ul	0.00
58) Fluorene-d10	15.73	176	77709	36.34	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.83	200	15113	32.00	ng/ul	0.00
71) Anthracene-d10	17.57	188	120522	36.82	ng/ul	0.00
79) Pyrene-d10	19.85	212	135193	37.00	ng/ul	0.00
90) Benzo(a)pyrene-d12	24.78	264	132871	36.07	ng/ul	0.01

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

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