

Data Path : Z:\HPCHEM1\BNA M\DATA\BM010116\
 Data File : BM003882.D
 Acq On : 30 Dec 2015 16:30
 Operator : SJ/UM
 Sample : PB87432BL
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK32

Quant Time: Dec 30 22:59:58 2015
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-BM010116.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Dec 30 22:56:39 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.67	152	49539	20.00	ng/ul	0.00
18) Naphthalene-d8	10.46	136	218833	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.33	164	123147	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.07	188	254910	20.00	ng/ul	0.00
78) Chrysene-d12	21.28	240	235265	20.00	ng/ul	0.00
86) Perylene-d12	23.51	264	223611	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.13	96	6879	6.21	ng/uL	0.00
5) Phenol-d5	6.85	99	125891	30.61	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.01	67	77224	33.22	ng/ul	0.00
9) 2-Chlorophenol-d4	7.20	132	107555	31.82	ng/ul	0.00
13) 4-Methylphenol-d8	8.39	113	104537	31.37	ng/ul	0.00
19) Nitrobenzene-d5	8.83	128	52548	31.88	ng/ul	0.00
22) 2-Nitrophenol-d4	9.55	143	57949	32.07	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.09	165	96556	30.08	ng/ul	0.00
29) 4-Chloroaniline-d4	10.60	131	152856	36.06	ng/ul	0.00
44) Dimethylphthalate-d6	13.74	166	318266	32.92	ng/ul	0.00
47) Acenaphthylene-d8	14.02	160	384328	32.63	ng/ul	0.00
52) 4-Nitrophenol-d4	14.54	143	48612	28.00	ng/ul	0.00
58) Fluorene-d10	15.32	176	268249	32.71	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.45	200	35941	25.00	ng/ul	0.00
71) Anthracene-d10	17.17	188	391263	33.19	ng/ul	0.00
79) Pyrene-d10	19.48	212	396921	32.58	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.37	264	377011	33.76	ng/ul	0.00

Target Compounds Ovalue

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

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