

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM011516\
 Data File : BM004017.D
 Acq On : 14 Jan 2016 17:56
 Operator : SJ/UM
 Sample : PB87811BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK11

Quant Time: Jan 15 00:31:39 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM010116.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Jan 15 00:28:21 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.66	152	37809	20.00	ng/ul	0.00
18) Naphthalene-d8	10.45	136	183748	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.32	164	108405	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.07	188	240997	20.00	ng/ul	0.00
78) Chrysene-d12	21.28	240	207281	20.00	ng/ul	0.00
86) Perylene-d12	23.52	264	181182	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.13	96	3833	4.53	ng/uL	0.00
5) Phenol-d5	6.84	99	86274	27.48	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.00	67	50091	28.23	ng/ul	0.00
9) 2-Chlorophenol-d4	7.20	132	72124	27.96	ng/ul	0.00
13) 4-Methylphenol-d8	8.38	113	74534	29.30	ng/ul	0.00
19) Nitrobenzene-d5	8.82	128	36895	26.66	ng/ul	0.00
22) 2-Nitrophenol-d4	9.55	143	40941	26.99	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.08	165	71145	26.39	ng/ul	0.00
29) 4-Chloroaniline-d4	10.59	131	112203	31.53	ng/ul	0.00
44) Dimethylphthalate-d6	13.73	166	241625	28.39	ng/ul	0.00
47) Acenaphthylene-d8	14.02	160	296376	28.59	ng/ul	0.00
52) 4-Nitrophenol-d4	14.54	143	34106	22.31	ng/ul	0.00
58) Fluorene-d10	15.32	176	209139	28.97	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.45	200	24008	17.67	ng/ul	0.00
71) Anthracene-d10	17.17	188	325849	29.23	ng/ul	0.00
79) Pyrene-d10	19.47	212	329654	30.71	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.37	264	269239	29.75	ng/ul	0.00

Target Compounds Qvalue

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

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