

Data Path : Z:\HPCHEM1\BNA M\DATA\BM100115\
 Data File : BM002810.D
 Acq On : 01 Oct 2015 18:47
 Operator : UM/IZ
 Sample : PB85865BL
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :

Quant Time: Oct 02 01:22:49 2015
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-BM092915.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Oct 02 01:21:34 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.69	152	66580	20.00	ng/ul	0.00
18) Naphthalene-d8	11.55	136	280460	20.00	ng/ul	0.00
36) Acenaphthene-d10	15.28	164	143888	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.99	188	302135	20.00	ng/ul	0.00
78) Chrysene-d12	22.08	240	327726	20.00	ng/ul	0.00
86) Perylene-d12	24.67	264	329268	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.76	96	7605	5.30	ng/uL	0.00
5) Phenol-d5	7.80	99	134942	26.75	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.98	67	76357	26.91	ng/ul	0.00
9) 2-Chlorophenol-d4	8.20	132	126025	26.84	ng/ul	0.00
13) 4-Methylphenol-d8	9.39	113	111541	26.82	ng/ul	0.00
19) Nitrobenzene-d5	9.88	128	58036	27.13	ng/ul	0.00
22) 2-Nitrophenol-d4	10.61	143	61307	27.43	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	11.15	165	99863	25.32	ng/ul	0.00
29) 4-Chloroaniline-d4	11.67	131	138138	29.09	ng/ul	0.00
44) Dimethylphthalate-d6	14.66	166	307422	28.40	ng/ul	0.00
47) Acenaphthylene-d8	14.99	160	391121	27.07	ng/ul	0.00
52) 4-Nitrophenol-d4	15.46	143	50712	26.38	ng/ul	0.00
58) Fluorene-d10	16.26	176	259824	26.64	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	16.38	200	31517	21.39	ng/ul	0.00
71) Anthracene-d10	18.09	188	397342	27.56	ng/ul	0.00
79) Pyrene-d10	20.33	212	416268	27.21	ng/ul	0.00
90) Benzo(a)pyrene-d12	24.50	264	471480	27.99	ng/ul	0.00

Target Compounds	Ovalue

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

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