

Data Path : Z:\HPCHEM1\BNA M\DATA\BM042516\
 Data File : BM005057.D
 Acq On : 25 Apr 2016 22:15
 Operator : UM/SJ
 Sample : PB90026BL
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK26

Quant Time: Apr 26 10:23:19 2016
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-BM042516.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Apr 26 00:45:11 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.80	152	30464	20.00	ng/ul	0.00
18) Naphthalene-d8	10.59	136	156607	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.44	164	121461	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.19	188	303023	20.00	ng/ul	0.00
75) Chrysene-d12	21.37	240	364406	20.00	ng/ul	0.00
83) Perylene-d12	23.66	264	327473	20.00	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	3.30	96	7409	12.36	ng/uL	0.00
5) Phenol-d5	6.96	99	140616	55.80	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.13	67	87274	60.87	ng/ul	0.00
9) 2-Chlorophenol-d4	7.33	132	115012	57.68	ng/ul	0.00
13) 4-Methylphenol-d8	8.50	113	120096	56.21	ng/ul	0.00
19) Nitrobenzene-d5	8.96	128	60203	55.50	ng/ul	0.00
22) 2-Nitrophenol-d4	9.67	143	69725	57.62	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.21	165	117496	51.11	ng/ul	0.00
29) 4-Chloroaniline-d4	10.73	131	180511	66.99	ng/ul	0.00
43) Dimethylphthalate-d6	13.85	166	453325	47.26	ng/ul	0.00
46) Acenaphthylene-d8	14.13	160	514352	45.03	ng/ul	0.00
51) 4-Nitrophenol-d4	14.64	143	70809	42.23	ng/ul	0.00
57) Fluorene-d10	15.43	176	377047	44.73	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.56	200	68867	39.96	ng/ul	0.00
70) Anthracene-d10	17.29	188	614145	45.20	ng/ul	0.00
76) Pyrene-d10	19.58	212	714965	43.10	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.52	264	684752	46.67	ng/ul	0.00

Target Compounds	Ovalue

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

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