

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM041415\
 Data File : BM000960.D
 Acq On : 13 Apr 2015 21:58
 Operator : TP/IZ
 Sample : PB82670BL
 Misc : SBLK43
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK43

Quant Time: Apr 14 18:51:31 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM01.2-EPA-BM041415.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Apr 14 08:28:33 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.60	152	139816	20.00	ng/ul	0.00
18) Naphthalene-d8	10.39	136	562859	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.26	164	310027	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.00	188	712853	20.00	ng/ul	0.00
75) Chrysene-d12	21.20	240	789918	20.00	ng/ul	0.00
83) Perylene-d12	23.38	264	746012	20.00	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	0.00	96	0d	0.00	ng/uL	
5) Phenol-d5	6.78	99	421582	40.11	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.94	67	252326	38.45	ng/ul	0.00
9) 2-Chlorophenol-d4	7.14	132	347903	40.10	ng/ul	0.00
13) 4-Methylphenol-d8	8.32	113	320433	39.13	ng/ul	0.00
19) Nitrobenzene-d5	8.76	128	142768	38.32	ng/ul	0.00
22) 2-Nitrophenol-d4	9.47	143	155412	38.47	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.01	165	293278	36.42	ng/ul	0.00
29) 4-Chloroaniline-d4	10.53	131	401644	41.64	ng/ul	0.00
43) Dimethylphthalate-d6	13.67	166	863345	42.53	ng/ul	0.00
46) Acenaphthylene-d8	13.95	160	1097143	37.35	ng/ul	0.00
51) 4-Nitrophenol-d4	14.47	143	104618	29.40	ng/ul	0.00
57) Fluorene-d10	15.26	176	789449	38.30	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.39	200	78680	20.66	ng/ul	0.00
70) Anthracene-d10	17.10	188	1247198	38.30	ng/ul	0.00
76) Pyrene-d10	19.40	212	1415670	39.70	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.24	264	1280973	39.25	ng/ul	0.00

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

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