

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM051215\
 Data File : BM001254.D
 Acq On : 12 May 2015 11:40
 Operator : TP/IZ
 Sample : PB83268BL
 Misc : SBLK57
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK57

Quant Time: May 12 18:39:31 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM051215.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue May 12 06:50:27 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.45	152	163122	20.00	ng/ul	0.00
18) Naphthalene-d8	10.23	136	688784	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.12	164	399142	20.00	ng/ul	0.00
61) Phenanthrene-d10	16.87	188	839239	20.00	ng/ul	0.00
75) Chrysene-d12	21.07	240	835300	20.00	ng/ul	0.00
83) Perylene-d12	23.19	264	787194	20.00	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	2.88	96	22174	6.62	ng/uL	0.00
5) Phenol-d5	6.63	99	471085	37.03	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.79	67	310861	39.65	ng/ul	0.00
9) 2-Chlorophenol-d4	6.97	132	393770	38.86	ng/ul	0.00
13) 4-Methylphenol-d8	8.17	113	409720	38.91	ng/ul	0.00
19) Nitrobenzene-d5	8.60	128	192512	42.38	ng/ul	0.00
22) 2-Nitrophenol-d4	9.32	143	216229	42.27	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	9.86	165	394622	38.81	ng/ul	0.00
29) 4-Chloroaniline-d4	10.37	131	598255	46.40	ng/ul	0.00
43) Dimethylphthalate-d6	13.54	166	1292055	47.76	ng/ul	0.00
46) Acenaphthylene-d8	13.81	160	1660510	42.84	ng/ul	0.00
51) 4-Nitrophenol-d4	14.35	143	195699	35.75	ng/ul	0.00
57) Fluorene-d10	15.12	176	1175034	43.47	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.26	200	201781	39.37	ng/ul	0.00
70) Anthracene-d10	16.97	188	1774669	46.34	ng/ul	0.00
76) Pyrene-d10	19.27	212	1871205	48.98	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.06	264	1664767	47.91	ng/ul	0.00

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

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