

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM042915\
 Data File : BM001133.D
 Acq On : 28 Apr 2015 18:09
 Operator : TP/IZ
 Sample : PB82970BL
 Misc : SBLK56
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK56

Quant Time: Apr 29 05:22:28 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM042815.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Apr 29 05:05:16 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.54	152	151112	20.00	ng/ul	0.00
18) Naphthalene-d8	10.32	136	637082	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.20	164	363483	20.00	ng/ul	0.00
61) Phenanthrene-d10	16.96	188	820620	20.00	ng/ul	0.00
75) Chrysene-d12	21.16	240	918645	20.00	ng/ul	0.00
83) Perylene-d12	23.32	264	918228	20.00	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	2.99	96	21018	6.30	ng/uL	0.00
5) Phenol-d5	6.72	99	443869	35.59	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.88	67	294848	37.26	ng/ul	0.00
9) 2-Chlorophenol-d4	7.07	132	356943	36.84	ng/ul	0.00
13) 4-Methylphenol-d8	8.26	113	352465	34.80	ng/ul	0.00
19) Nitrobenzene-d5	8.69	128	163308	37.94	ng/ul	0.00
22) 2-Nitrophenol-d4	9.42	143	180051	38.50	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	9.95	165	319817	34.66	ng/ul	0.00
29) 4-Chloroaniline-d4	10.47	131	496561	45.56	ng/ul	0.00
43) Dimethylphthalate-d6	13.62	166	1050160	41.79	ng/ul	0.00
46) Acenaphthylene-d8	13.90	160	1326183	37.33	ng/ul	0.00
51) 4-Nitrophenol-d4	14.42	143	156234	29.55	ng/ul	0.00
57) Fluorene-d10	15.20	176	944265	38.02	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.33	200	150036	30.52	ng/ul	0.00
70) Anthracene-d10	17.06	188	1467819	38.60	ng/ul	0.00
76) Pyrene-d10	19.36	212	1670846	41.03	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.19	264	1649125	39.61	ng/ul	0.00

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

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