

Data Path : Z:\HPCHEM1\BNA M\DATA\BM081716\
 Data File : BM007143.D
 Acq On : 16 Aug 2016 16:01
 Operator : UM/SJ
 Sample : PB92870BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK70

Quant Time: Aug 17 02:28:49 2016
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-BM081116.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Aug 17 02:24:57 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.81	152	213787	20.00	ng/ul	0.00
18) Naphthalene-d8	10.59	136	857810	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.44	164	512940	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.18	188	1211816	20.00	ng/ul	0.00
75) Chrysene-d12	21.36	240	1549844	20.00	ng/ul	0.00
83) Perylene-d12	23.64	264	1612051	20.00	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	3.33	96	36322	7.42	ng/uL	0.00
5) Phenol-d5	6.97	99	621713	35.90	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.15	67	368980	37.17	ng/ul	0.00
9) 2-Chlorophenol-d4	7.34	132	511404	37.07	ng/ul	0.00
13) 4-Methylphenol-d8	8.50	113	504131	35.37	ng/ul	0.00
19) Nitrobenzene-d5	8.96	128	236435	37.95	ng/ul	0.00
22) 2-Nitrophenol-d4	9.68	143	264661	39.76	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.21	165	495648	35.55	ng/ul	0.00
29) 4-Chloroaniline-d4	10.73	131	658512	39.69	ng/ul	0.00
43) Dimethylphthalate-d6	13.85	166	1577621	37.22	ng/ul	0.00
46) Acenaphthylene-d8	14.13	160	1892997	37.18	ng/ul	0.00
51) 4-Nitrophenol-d4	14.63	143	233349	30.72	ng/ul	0.00
57) Fluorene-d10	15.43	176	1332866	35.98	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.54	200	197977	29.56	ng/ul	0.00
70) Anthracene-d10	17.28	188	2146750	38.24	ng/ul	0.00
76) Pyrene-d10	19.57	212	2492911	38.11	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.50	264	2810714	38.12	ng/ul	0.00

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

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