

Data Path : Z:\HPCHEM1\BNA M\DATA\BM081816\
 Data File : BM007184.D
 Acq On : 17 Aug 2016 18:09
 Operator : UM/SJ
 Sample : PB92841BL
 Misc :
 ALS Vial : 39 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK41

Quant Time: Aug 17 18:40:21 2016
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-BM081116.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Aug 17 15:21:18 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.80	152	250237	20.00	ng/ul	0.00
18) Naphthalene-d8	10.59	136	999975	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.43	164	581731	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.18	188	1368866	20.00	ng/ul	0.00
75) Chrysene-d12	21.36	240	1774729	20.00	ng/ul	0.00
83) Perylene-d12	23.63	264	1845748	20.00	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	3.32	96	41869	7.31	ng/uL	0.00
5) Phenol-d5	6.97	99	736019	36.31	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.14	67	428644	36.89	ng/ul	0.00
9) 2-Chlorophenol-d4	7.34	132	602120	37.29	ng/ul	0.00
13) 4-Methylphenol-d8	8.50	113	585317	35.09	ng/ul	0.00
19) Nitrobenzene-d5	8.96	128	279424	38.47	ng/ul	0.00
22) 2-Nitrophenol-d4	9.67	143	311480	40.14	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.21	165	585126	36.00	ng/ul	0.00
29) 4-Chloroaniline-d4	10.73	131	752451	38.91	ng/ul	0.00
43) Dimethylphthalate-d6	13.84	166	1752915	36.47	ng/ul	0.00
46) Acenaphthylene-d8	14.13	160	2138976	37.05	ng/ul	0.00
51) 4-Nitrophenol-d4	14.63	143	269338	31.26	ng/ul	0.00
57) Fluorene-d10	15.43	176	1502771	35.77	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.54	200	209065	27.64	ng/ul	0.00
70) Anthracene-d10	17.27	188	2394097	37.75	ng/ul	0.00
76) Pyrene-d10	19.57	212	2839832	37.91	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.49	264	3256042	38.57	ng/ul	0.00

Target Compounds	Ovalue

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

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