

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM111116\
 Data File : BM007832.D
 Acq On : 11 Nov 2016 18:59
 Operator : UM/SJ
 Sample : PB94658BL
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK58

Quant Time: Nov 12 03:40:45 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM102016.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Nov 12 03:12:55 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.74	152	164597	20.00	ng/ul	0.00
18) Naphthalene-d8	10.52	136	639253	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.37	164	409110	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.12	188	811816	20.00	ng/ul	0.00
75) Chrysene-d12	21.31	240	1003351	20.00	ng/ul	0.00
83) Perylene-d12	23.56	264	1039671	20.00	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	3.28	96	28248	7.35	ng/uL	0.00
5) Phenol-d5	6.92	99	481103	37.87	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.07	67	287946	38.19	ng/ul	0.00
9) 2-Chlorophenol-d4	7.27	132	384319	39.45	ng/ul	0.00
13) 4-Methylphenol-d8	8.45	113	379432	38.17	ng/ul	0.00
19) Nitrobenzene-d5	8.89	128	180468	39.23	ng/ul	0.00
22) 2-Nitrophenol-d4	9.61	143	198735	40.75	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.15	165	356664	37.20	ng/ul	0.00
29) 4-Chloroaniline-d4	10.66	131	471548	43.11	ng/ul	0.00
43) Dimethylphthalate-d6	13.79	166	1117314	40.11	ng/ul	0.00
46) Acenaphthylene-d8	14.06	160	1338041	39.91	ng/ul	0.00
51) 4-Nitrophenol-d4	14.59	143	146273	33.21	ng/ul	0.00
57) Fluorene-d10	15.37	176	962622	39.43	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.51	200	127385	26.41	ng/ul	0.00
70) Anthracene-d10	17.22	188	1523243	41.64	ng/ul	0.00
76) Pyrene-d10	19.52	212	1761560	42.73	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.42	264	1913546	42.16	ng/ul	0.00

Target Compounds	Qvalue

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

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