

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM121416\
 Data File : BM008310.D
 Acq On : 14 Dec 2016 15:58
 Operator : UM/SJ
 Sample : PB95250BL
 Misc :
 ALS Vial : 36 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK50

Quant Time: Dec 14 16:58:28 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-2016\SOM02.2-EPA-BM121416.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Dec 14 05:32:31 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.67	152	111853	20.00	ng/ul	0.00
18) Naphthalene-d8	10.45	136	441346	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.31	164	288173	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.06	188	582077	20.00	ng/ul	0.00
75) Chrysene-d12	21.27	240	732597	20.00	ng/ul	0.00
83) Perylene-d12	23.50	264	758288	20.00	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	3.24	96	19467	8.19	ng/uL	0.00
5) Phenol-d5	6.86	99	308823	35.59	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.02	67	189242	37.94	ng/ul	0.00
9) 2-Chlorophenol-d4	7.21	132	249916	38.28	ng/ul	0.00
13) 4-Methylphenol-d8	8.39	113	244401	36.36	ng/ul	0.00
19) Nitrobenzene-d5	8.83	128	122376	38.37	ng/ul	0.00
22) 2-Nitrophenol-d4	9.55	143	136311	38.57	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.09	165	225177	34.76	ng/ul	0.00
29) 4-Chloroaniline-d4	10.61	131	259216	34.91	ng/ul	0.00
43) Dimethylphthalate-d6	13.73	166	756163	39.47	ng/ul	0.00
46) Acenaphthylene-d8	14.00	160	898778	39.70	ng/ul	0.00
51) 4-Nitrophenol-d4	14.57	143	83137	29.46	ng/ul	0.01
57) Fluorene-d10	15.31	176	640913	38.49	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.47	200	93645	27.03	ng/ul	0.00
70) Anthracene-d10	17.16	188	1006041	39.67	ng/ul	0.00
76) Pyrene-d10	19.47	212	1176895	41.60	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.35	264	1260718	39.59	ng/ul	0.00

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

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