

Data Path : Z:\HPCHEM1\BNA M\DATA\BM102016\
 Data File : BM007707.D
 Acq On : 20 Oct 2016 17:29
 Operator : UM/SJ
 Sample : PB94208BL
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK08

Quant Time: Oct 21 10:08:57 2016
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-BM102016.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Oct 20 17:01:48 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.76	152	162597	20.00	ng/ul	-0.02
18) Naphthalene-d8	10.53	136	611689	20.00	ng/ul	-0.02
35) Acenaphthene-d10	14.39	164	385480	20.00	ng/ul	-0.02
61) Phenanthrene-d10	17.13	188	765561	20.00	ng/ul	-0.02
75) Chrysene-d12	21.32	240	973766	20.00	ng/ul	-0.02
83) Perylene-d12	23.57	264	1002083	20.00	ng/ul	-0.02

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	3.29	96	27663	7.29	ng/uL	0.00
5) Phenol-d5	6.93	99	456103	36.34	ng/ul	-0.02
7) Bis-(2-Chloroethyl)ether-d	7.09	67	278799	37.44	ng/ul	-0.02
9) 2-Chlorophenol-d4	7.29	132	370353	38.48	ng/ul	-0.02
13) 4-Methylphenol-d8	8.46	113	360937	36.76	ng/ul	-0.02
19) Nitrobenzene-d5	8.91	128	172815	39.26	ng/ul	-0.02
22) 2-Nitrophenol-d4	9.63	143	187947	40.27	ng/ul	-0.02
26) 2,4-Dichlorophenol-d3	10.16	165	340901	37.16	ng/ul	-0.02
29) 4-Chloroaniline-d4	10.67	131	454226	43.40	ng/ul	-0.02
43) Dimethylphthalate-d6	13.80	166	1096024	41.76	ng/ul	-0.02
46) Acenaphthylene-d8	14.08	160	1265052	40.05	ng/ul	-0.02
51) 4-Nitrophenol-d4	14.60	143	136731	32.95	ng/ul	-0.02
57) Fluorene-d10	15.38	176	914932	39.77	ng/ul	-0.02
62) 4,6-Dinitro-2-methylphenol	15.52	200	145433	31.97	ng/ul	-0.01
70) Anthracene-d10	17.23	188	1441260	41.78	ng/ul	-0.02
76) Pyrene-d10	19.53	212	1664480	41.60	ng/ul	-0.02
87) Benzo(a)pyrene-d12	23.43	264	1842525	42.12	ng/ul	-0.02

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

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