

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM011616\
 Data File : BM004042.D
 Acq On : 15 Jan 2016 11:46
 Operator : SJ/UM
 Sample : PB87822BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK22

Quant Time: Jan 16 00:54:06 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM010116.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Jan 16 00:22:06 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.66	152	35977	20.00	ng/ul	0.00
18) Naphthalene-d8	10.45	136	175233	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.32	164	101458	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.07	188	214773	20.00	ng/ul	0.00
78) Chrysene-d12	21.27	240	170201	20.00	ng/ul	0.00
86) Perylene-d12	23.51	264	159778	20.00	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	3.13	96	5044	6.27	ng/uL	0.00
5) Phenol-d5	6.84	99	112079	37.52	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.00	67	65587	38.84	ng/ul	0.00
9) 2-Chlorophenol-d4	7.20	132	93899	38.25	ng/ul	0.00
13) 4-Methylphenol-d8	8.37	113	97063	40.10	ng/ul	0.00
19) Nitrobenzene-d5	8.82	128	48981	37.11	ng/ul	0.00
22) 2-Nitrophenol-d4	9.55	143	54538	37.69	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.08	165	92399	35.94	ng/ul	0.00
29) 4-Chloroaniline-d4	10.59	131	145442	42.85	ng/ul	0.00
44) Dimethylphthalate-d6	13.73	166	307312	38.58	ng/ul	0.00
47) Acenaphthylene-d8	14.01	160	380175	39.18	ng/ul	0.00
52) 4-Nitrophenol-d4	14.54	143	42350	29.60	ng/ul	0.00
58) Fluorene-d10	15.32	176	267250	39.56	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.45	200	27139	22.41	ng/ul	0.00
71) Anthracene-d10	17.17	188	398651	40.13	ng/ul	0.00
79) Pyrene-d10	19.47	212	370614	42.04	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.37	264	329624	41.31	ng/ul	0.00

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

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