

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM040616\
 Data File : BM004903.D
 Acq On : 06 Apr 2016 14:00
 Operator : UM/SJ
 Sample : PB89199BL
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK99

Quant Time: Apr 07 04:32:28 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM040116.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Apr 07 04:29:45 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.82	152	55614	20.00	ng/ul	0.00
18) Naphthalene-d8	10.61	136	251698	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.46	164	164453	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.20	188	416950	20.00	ng/ul	0.00
78) Chrysene-d12	21.39	240	559038	20.00	ng/ul	0.00
86) Perylene-d12	23.68	264	556881	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	0.00	96	0	0.00	ng/uL	
5) Phenol-d5	6.99	99	1686	0.34	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.16	67	1523	0.57	ng/ul	0.00
9) 2-Chlorophenol-d4	7.36	132	1603	0.43	ng/ul	0.00
13) 4-Methylphenol-d8	8.53	113	1082	0.26	ng/ul	0.02
19) Nitrobenzene-d5	8.98	128	361	0.20	ng/ul	0.00
22) 2-Nitrophenol-d4	9.70	143	388	0.20	ng/ul	0.01
26) 2,4-Dichlorophenol-d3	10.25	165	683	0.18	ng/ul	0.02
29) 4-Chloroaniline-d4	10.76	131	6450	1.46	ng/ul	0.01
44) Dimethylphthalate-d6	13.86	166	7622	0.59	ng/ul	0.00
47) Acenaphthylene-d8	14.15	160	9076	0.58	ng/ul	0.00
52) 4-Nitrophenol-d4	0.00	143	0	0.00	ng/ul	
58) Fluorene-d10	15.45	176	9029	0.81	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	0.00	200	0	0.00	ng/ul	
71) Anthracene-d10	17.30	188	15238	0.82	ng/ul	0.00
79) Pyrene-d10	19.59	212	16852	0.69	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.53	264	19416	0.79	ng/ul	0.00

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

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