

Data Path : Z:\HPCHEM1\BNA M\DATA\BM071416\
 Data File : BM006418.D
 Acq On : 14 Jul 2016 19:17
 Operator : UM/SJ
 Sample : PB92088BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK88

Quant Time: Jul 15 04:40:49 2016
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-BM071416.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Jul 15 04:34:07 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.62	152	86854	20.00	ng/ul	0.00
18) Naphthalene-d8	10.40	136	364793	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.27	164	229099	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.02	188	566901	20.00	ng/ul	0.00
75) Chrysene-d12	21.22	240	756086	20.00	ng/ul	0.00
83) Perylene-d12	23.42	264	804514	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.14	96	15061	7.30	ng/uL	0.00
5) Phenol-d5	6.80	99	242126	34.02	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.96	67	152476	35.32	ng/ul	0.00
9) 2-Chlorophenol-d4	7.16	132	197936	34.32	ng/ul	0.00
13) 4-Methylphenol-d8	8.33	113	193041	34.13	ng/ul	0.00
19) Nitrobenzene-d5	8.77	128	93119	33.68	ng/ul	0.00
22) 2-Nitrophenol-d4	9.49	143	109577	35.05	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.03	165	188102	31.89	ng/ul	0.00
29) 4-Chloroaniline-d4	10.54	131	255716	41.48	ng/ul	0.00
43) Dimethylphthalate-d6	13.69	166	664900	35.46	ng/ul	0.00
46) Acenaphthylene-d8	13.96	160	773334	33.36	ng/ul	0.00
51) 4-Nitrophenol-d4	14.49	143	109480	33.61	ng/ul	0.00
57) Fluorene-d10	15.27	176	560866	33.65	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.40	200	95116	29.00	ng/ul	0.00
70) Anthracene-d10	17.12	188	925268	34.53	ng/ul	0.00
76) Pyrene-d10	19.42	212	1106444	32.28	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.28	264	1303482	35.43	ng/ul	0.00

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

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