

Data Path : Z:\HPCHEM1\BNA M\DATA\BM022616\
 Data File : BM004425.D
 Acq On : 26 Feb 2016 12:21
 Operator : SJ/UM
 Sample : H1572-01
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 0311-RIN07

Quant Time: Feb 26 23:25:10 2016
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-BM022216.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Feb 26 23:20:20 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.61	152	22871	20.00	ng/ul	0.00
18) Naphthalene-d8	10.40	136	103724	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.28	164	61948	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.03	188	128944	20.00	ng/ul	0.00
78) Chrysene-d12	21.25	240	125640	20.00	ng/ul	0.00
86) Perylene-d12	23.47	264	123065	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.08	96	507	1.17	ng/uL	0.00
5) Phenol-d5	6.81	99	12490	6.51	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.95	67	28821	29.55	ng/ul	0.00
9) 2-Chlorophenol-d4	7.15	132	40827	24.68	ng/ul	0.00
13) 4-Methylphenol-d8	8.34	113	25374	14.81	ng/ul	0.00
19) Nitrobenzene-d5	8.77	128	24268	30.75	ng/ul	0.00
22) 2-Nitrophenol-d4	9.50	143	27443	30.51	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.04	165	43561	26.79	ng/ul	0.00
29) 4-Chloroaniline-d4	0.00	131	0	0.00	ng/ul	
44) Dimethylphthalate-d6	13.69	166	165818	31.21	ng/ul	0.00
47) Acenaphthylene-d8	13.97	160	203532	32.72	ng/ul	0.00
52) 4-Nitrophenol-d4	14.56	143	2391	2.80	ng/ul	0.04
58) Fluorene-d10	15.28	176	142069	31.44	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.43	200	17412	21.98	ng/ul	0.00
71) Anthracene-d10	17.13	188	210195	33.27	ng/ul	0.00
79) Pyrene-d10	19.44	212	216686	35.65	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.33	264	218557	33.84	ng/ul	0.00

Target Compounds Ovalue

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

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