

Data Path : Z:\HPCHEM1\BNA M\DATA\BM030518\
 Data File : BM014488.D
 Acq On : 06 Mar 2018 00:23
 Operator : SJ/JU
 Sample : J1795-10
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :

Quant Time: Mar 06 03:33:15 2018
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM021418.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Mar 05 23:57:03 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.51	152	71725	20.00	ng/ul	0.00
18) Naphthalene-d8	10.29	136	260545	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.17	164	148981	20.00	ng/ul	0.00
61) Phenanthrene-d10	16.94	188	322624	20.00	ng/ul	0.00
75) Chrysene-d12	21.16	240	300795	20.00	ng/ul	0.00
83) Perylene-d12	23.34	264	294691	20.00	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	3.06	96	221	0.13	ng/uL	-0.03
5) Phenol-d5	0.00	99	0	0.00	ng/ul	
7) Bis-(2-Chloroethyl)ether-d	0.00	67	0	0.00	ng/ul	
9) 2-Chlorophenol-d4	0.00	132	0	0.00	ng/ul	
13) 4-Methylphenol-d8	0.00	113	0	0.00	ng/ul	
19) Nitrobenzene-d5	0.00	128	0	0.00	ng/ul	
22) 2-Nitrophenol-d4	0.00	143	0	0.00	ng/ul	
26) 2,4-Dichlorophenol-d3	0.00	165	0	0.00	ng/ul	
29) 4-Chloroaniline-d4	0.00	131	0	0.00	ng/ul	
43) Dimethylphthalate-d6	0.00	166	0	0.00	ng/ul	
46) Acenaphthylene-d8	0.00	160	0	0.00	ng/ul	
51) 4-Nitrophenol-d4	0.00	143	0	0.00	ng/ul	
57) Fluorene-d10	15.21	176	256	0.02	ng/ul	0.03
62) 4,6-Dinitro-2-methylphenol	0.00	200	0	0.00	ng/ul	
70) Anthracene-d10	16.94	188	322624	20.73	ng/ul	-0.09
76) Pyrene-d10	19.46	212	303	0.02	ng/ul	0.11
87) Benzo(a)pyrene-d12	23.34	264	289667	20.20	ng/ul	0.13

Target Compounds	Ovalue

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

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