

Data Path : Z:\HPCHEM1\BNA M\DATA\BM073016\
 Data File : BM006775.D
 Acq On : 30 Jul 2016 16:10
 Operator : UM/SJ
 Sample : H4190-14
 Misc :
 ALS Vial : 38 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :

Quant Time: Aug 01 04:25:55 2016
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-BM072016.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Jul 30 01:37:26 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.60	152	171291	20.00	ng/ul	0.00
7) Naphthalene-d8	10.38	136	777014	20.00	ng/ul	0.00
15) Acenaphthene-d10	14.26	164	449103	20.00	ng/ul	0.00
23) Phenanthrene-d10	17.01	188	1005954	20.00	ng/ul	0.00
29) Chrysene-d12	21.21	240	1132069	20.00	ng/ul	0.00
34) Perylene-d12	23.41	264	1167526	20.00	ng/ul	0.00

System Monitoring Compounds

2) 1,4-Dioxane-d8	3.13	96	8113	1.92	ng/uL	0.00
3) Phenol-d5	6.79	99	298989	20.52	ng/ul	0.00
4) Bis-(2-Chloroethyl)ether-d	6.95	67	211852	23.47	ng/ul	0.00
5) 2-Chlorophenol-d4	7.14	132	247977	21.30	ng/ul	0.00
6) 4-Methylphenol-d8	8.32	113	169485	15.00	ng/ul	0.00
8) Nitrobenzene-d5	8.76	128	127634	21.70	ng/ul	0.00
9) 2-Nitrophenol-d4	9.47	143	141479	22.62	ng/ul	0.00
10) 2,4-Dichlorophenol-d3	10.01	165	249038	21.07	ng/ul	0.00
12) 4-Chloroaniline-d4	10.53	131	334639	25.71	ng/ul	0.00
16) Dimethylphthalate-d6	13.67	166	838635	23.82	ng/ul	0.00
17) Acenaphthylene-d8	13.94	160	1065952	23.72	ng/ul	0.00
20) 4-Nitrophenol-d4	14.49	143	122300	20.00	ng/ul	0.00
21) Fluorene-d10	15.26	176	744298	23.63	ng/ul	0.00
24) 4,6-Dinitro-2-methylphenol	15.40	200	97768	18.72	ng/ul	0.00
26) Anthracene-d10	17.10	188	1131027	23.68	ng/ul	0.00
30) Pyrene-d10	19.41	212	1306430	25.33	ng/ul	0.00
37) Benzo(a)pyrene-d12	23.27	264	1325705	24.69	ng/ul	0.00

Target Compounds	Ovalue

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

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