

Data Path : Z:\HPCHEM1\BNA M\DATA\BM073016\
 Data File : BM006783.D
 Acq On : 31 Jul 2016 11:17
 Operator : UM/SJ
 Sample : H4188-13
 Misc :
 ALS Vial : 52 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :

Quant Time: Aug 01 04:26:56 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	210068	20.00	ng/ul	0.00
7) Naphthalene-d8	10.38	136	901787	20.00	ng/ul	0.00
15) Acenaphthene-d10	14.26	164	498450	20.00	ng/ul	0.00
23) Phenanthrene-d10	17.01	188	1113644	20.00	ng/ul	0.00
29) Chrysene-d12	21.21	240	1282384	20.00	ng/ul	0.00
34) Perylene-d12	23.41	264	1317025	20.00	ng/ul	0.00

System Monitoring Compounds

2) 1,4-Dioxane-d8	3.13	96	8250	1.59	ng/uL	0.00
3) Phenol-d5	6.79	99	330005	18.47	ng/ul	0.00
4) Bis-(2-Chloroethyl)ether-d	6.95	67	216790	19.58	ng/ul	0.00
5) 2-Chlorophenol-d4	7.14	132	264431	18.52	ng/ul	0.00
6) 4-Methylphenol-d8	8.32	113	235378	16.98	ng/ul	0.00
8) Nitrobenzene-d5	8.76	128	130093	19.06	ng/ul	0.00
9) 2-Nitrophenol-d4	9.47	143	144112	19.85	ng/ul	0.00
10) 2,4-Dichlorophenol-d3	10.01	165	253007	18.44	ng/ul	0.00
12) 4-Chloroaniline-d4	10.53	131	330798	21.90	ng/ul	0.00
16) Dimethylphthalate-d6	13.67	166	786297	20.12	ng/ul	0.00
17) Acenaphthylene-d8	13.95	160	1009384	20.23	ng/ul	0.00
20) 4-Nitrophenol-d4	14.49	143	105379	15.52	ng/ul	0.00
21) Fluorene-d10	15.25	176	708122	20.26	ng/ul	0.00
24) 4,6-Dinitro-2-methylphenol	15.40	200	95453	16.51	ng/ul	0.00
26) Anthracene-d10	17.10	188	1057737	20.00	ng/ul	0.00
30) Pyrene-d10	19.41	212	1241458	21.25	ng/ul	0.00
37) Benzo(a)pyrene-d12	23.27	264	1265216	20.88	ng/ul	-0.01

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

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