

Data Path : Z:\HPCHEM1\BNA M\DATA\BM102517\
 Data File : BM012473.D
 Acq On : 26 Oct 2017 15:26
 Operator : SJ/JU
 Sample : I5882-01DL 50X
 Misc :
 ALS Vial : 36 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DAHH0DL

Quant Time: Oct 26 17:42:55 2017
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM102417.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Oct 26 09:35:31 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.88	152	539416	20.00	ng/ul	0.00
18) Naphthalene-d8	10.67	136	2218578	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.50	164	1308663	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.25	188	2622738	20.00	ng/ul	0.00
75) Chrysene-d12	21.42	240	2124159	20.00	ng/ul	0.00
83) Perylene-d12	23.72	264	2372518	20.00	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.37	96	1747	0.17	ng/uL	0.00
5) Phenol-d5	7.05	99	18253	0.41	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.21	67	14767	0.55	ng/ul	0.00
9) 2-Chlorophenol-d4	7.42	132	18085	0.54	ng/ul	0.00
13) 4-Methylphenol-d8	8.58	113	16269	0.42	ng/ul	0.00
19) Nitrobenzene-d5	9.03	128	9510	0.68	ng/ul	0.00
22) 2-Nitrophenol-d4	9.75	143	9219	0.64	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.29	165	19624	0.52	ng/ul	-0.01
29) 4-Chloroaniline-d4	10.80	131	18078	0.45	ng/ul	0.00
43) Dimethylphthalate-d6	13.92	166	64413	0.56	ng/ul	0.00
46) Acenaphthylene-d8	14.20	160	74194	0.55	ng/ul	0.00
51) 4-Nitrophenol-d4	14.70	143	6603	0.39	ng/ul	0.00
57) Fluorene-d10	15.50	176	53874	0.56	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.63	200	5261	0.41	ng/ul	0.01
70) Anthracene-d10	17.35	188	82538	0.66	ng/ul	0.00
76) Pyrene-d10	19.63	212	65296	0.67	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.57	264	68256	0.61	ng/ul	0.00
Target Compounds						
55) 2,3,4,6-Tetrachlorophenol	15.13	232	467972	17.190	ng/ul	96
68) Pentachlorophenol	16.90	266	5233100	275.317	ng/ul	97

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

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