

Data Path : Z:\HPCHEM1\BNA_M\Data\BM100317\
 Data File : BM011881.D
 Acq On : 04 Oct 2017 09:17
 Operator : SJ/JU
 Sample : I5354-10
 Misc :
 ALS Vial : 35 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 B0BA4

Quant Time: Oct 04 12:35:57 2017
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM100317.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Oct 04 05:30:09 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.87	152	212379	20.00	ng/ul	0.00
18) Naphthalene-d8	10.67	136	949675	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.52	164	624015	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.26	188	1536424	20.00	ng/ul	0.00
75) Chrysene-d12	21.43	240	1901353	20.00	ng/ul	0.00
83) Perylene-d12	23.75	264	1826574	20.00	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.29	96	25091	5.59	ng/uL	0.00
5) Phenol-d5	7.03	99	450230	24.64	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.20	67	277896	25.70	ng/ul	0.00
9) 2-Chlorophenol-d4	7.40	132	381559	26.03	ng/ul	0.00
13) 4-Methylphenol-d8	8.57	113	405111	25.65	ng/ul	0.00
19) Nitrobenzene-d5	9.04	128	183299	27.10	ng/ul	0.00
22) 2-Nitrophenol-d4	9.76	143	202436	26.57	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.30	165	412896	26.82	ng/ul	0.00
29) 4-Chloroaniline-d4	10.82	131	449435	26.76	ng/ul	0.00
43) Dimethylphthalate-d6	13.92	166	1445424	26.73	ng/ul	0.00
46) Acenaphthylene-d8	14.21	160	1716359	27.01	ng/ul	0.00
51) 4-Nitrophenol-d4	14.71	143	176568	18.69	ng/ul	0.00
57) Fluorene-d10	15.50	176	1256598	26.33	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.63	200	197945	19.15	ng/ul	0.00
70) Anthracene-d10	17.36	188	2058929	27.20	ng/ul	0.00
76) Pyrene-d10	19.64	212	2558397	29.83	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.60	264	2447928	27.86	ng/ul	0.00
Target Compounds						
4) Benzaldehyde	7.02	77	12808	1.407	ng/ul	95
6) Phenol	7.06	94	64387	3.561	ng/ul	89
14) Acetophenone	8.70	105	396800	17.168	ng/ul	91

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

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