

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM100418\
 Data File : BM017010.D
 Acq On : 04 Oct 2018 21:25
 Operator : MJ/SJ
 Sample : J5157-07
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0AC9

Quant Time: Oct 05 05:19:13 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM100418MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Oct 05 04:50:45 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.70	152	201631	20.00	ng/ul	0.00
18) Naphthalene-d8	10.47	136	879808	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.34	164	477425	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.09	188	1064343	20.00	ng/ul	0.00
78) Chrysene-d12	21.27	240	1145142	20.00	ng/ul	0.00
86) Perylene-d12	23.50	264	1186990	20.00	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.25	96	23717	5.85	ng/uL	0.00
5) Phenol-d5	6.87	99	527122	31.27	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.04	67	345652	33.82	ng/ul	0.00
9) 2-Chlorophenol-d4	7.23	132	455477	32.50	ng/ul	0.00
13) 4-Methylphenol-d8	8.40	113	437715	33.14	ng/ul	0.00
19) Nitrobenzene-d5	8.85	128	222915	34.67	ng/ul	0.00
22) 2-Nitrophenol-d4	9.57	143	230979	35.31	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.10	165	434594	32.29	ng/ul	0.00
29) 4-Chloroaniline-d4	10.62	131	630574	38.99	ng/ul	0.00
44) Dimethylphthalate-d6	13.75	166	1392495	36.76	ng/ul	0.00
47) Acenaphthylene-d8	14.03	160	1768979	36.54	ng/ul	0.00
52) 4-Nitrophenol-d4	14.54	143	168506	23.54	ng/ul	0.00
58) Fluorene-d10	15.33	176	1232849	36.54	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.46	200	128023	20.40	ng/ul	0.00
71) Anthracene-d10	17.19	188	1856369	36.52	ng/ul	0.00
79) Pyrene-d10	19.48	212	2023248	39.17	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.36	264	2300800	37.86	ng/ul	0.00
Target Compounds						
6) Phenol	6.90	94	28664	1.700	ng/ul#	88
45) Dimethylphthalate	13.80	163	624363	16.646	ng/ul	100

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

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