

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM120319\
 Data File : BM023971.D
 Acq On : 03 Dec 2019 14:05
 Operator : JU
 Sample : K5914-03
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 JLNR4

Quant Time: Dec 03 14:56:57 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM112719MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Dec 03 11:33:32 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.49	152	439927	20.00	ng/ul	0.00
18) Naphthalene-d8	10.26	136	1953709	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.14	164	1264714	20.00	ng/ul	0.00
62) Phenanthrene-d10	16.90	188	2623096	20.00	ng/ul	0.00
78) Chrysene-d12	21.11	240	1822100	20.00	ng/ul	0.00
86) Perylene-d12	23.25	264	1953541	20.00	ng/ul	0.00

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.04	96	39615	4.01	ng/uL	0.00
5) Phenol-d5	6.67	99	829665	21.52	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.83	67	515326	22.98	ng/ul	0.00
9) 2-Chlorophenol-d4	7.02	132	657584	22.07	ng/ul	0.00
13) 4-Methylphenol-d8	8.20	113	646066	21.24	ng/ul	0.00
19) Nitrobenzene-d5	8.63	128	337984	22.29	ng/ul	0.00
22) 2-Nitrophenol-d4	9.35	143	385372	22.36	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	9.89	165	684015	20.96	ng/ul	0.00
29) 4-Chloroaniline-d4	10.40	131	823664	23.00	ng/ul	0.00
44) Dimethylphthalate-d6	13.57	166	2318634	23.59	ng/ul	0.00
47) Acenaphthylene-d8	13.83	160	2732241	23.72	ng/ul	0.00
52) 4-Nitrophenol-d4	14.37	143	253973	15.37	ng/ul	0.00
58) Fluorene-d10	15.14	176	1996612	24.14	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.29	200	244612	14.85	ng/ul	0.00
71) Anthracene-d10	17.00	188	3049896	24.34	ng/ul	0.00
79) Pyrene-d10	19.30	212	3166375	33.63	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.11	264	2630578	25.47	ng/ul	0.00

Target Compounds					Ovalue	
6) Phenol	6.70	94	70266	1.772	ng/ul	95
45) Dimethylphthalate	13.62	163	938240	9.754	ng/ul	99

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

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