

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM100319\
 Data File : BM023011.D
 Acq On : 06 Oct 2019 04:05
 Operator : JU
 Sample : K5057-07
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DBAJ2

Quant Time: Oct 07 03:45:58 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM092719MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Oct 04 16:45:59 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.79	152	251963	20.00	ng/ul	0.00
18) Naphthalene-d8	10.57	136	1113919	20.00	ng/ul	-0.01
36) Acenaphthene-d10	14.42	164	678334	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.16	188	1216856	20.00	ng/ul	0.00
78) Chrysene-d12	21.34	240	903142	20.00	ng/ul	-0.01
86) Perylene-d12	23.60	264	989145	20.00	ng/ul	-0.01

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.27	96	14325	2.46	ng/uL	0.00
5) Phenol-d5	6.95	99	254037	11.41	ng/ul	-0.01
7) Bis-(2-Chloroethyl)ether-d	7.12	67	157644	12.94	ng/ul	0.00
9) 2-Chlorophenol-d4	7.32	132	211578	11.80	ng/ul	0.00
13) 4-Methylphenol-d8	8.49	113	214834	11.79	ng/ul	-0.01
19) Nitrobenzene-d5	8.94	128	106562	12.40	ng/ul	0.00
22) 2-Nitrophenol-d4	9.66	143	103551	11.03	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.20	165	203043	10.86	ng/ul	-0.01
29) 4-Chloroaniline-d4	10.71	131	288460	12.70	ng/ul	0.00
44) Dimethylphthalate-d6	13.83	166	700150	13.30	ng/ul	0.00
47) Acenaphthylene-d8	14.12	160	809242	13.31	ng/ul	0.00
52) 4-Nitrophenol-d4	14.62	143	23855	2.29	ng/ul	0.00
58) Fluorene-d10	15.42	176	596425	13.29	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.53	200	31460	3.97	ng/ul	-0.01
71) Anthracene-d10	17.26	188	866658	14.58	ng/ul	-0.01
79) Pyrene-d10	19.55	212	863229	17.76	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.45	264	746911	14.60	ng/ul	-0.02

Target Compounds					Ovalue	
6) Phenol	6.98	94	47930	2.054	ng/ul	97
45) Dimethylphthalate	13.88	163	233241	4.313	ng/ul	100

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

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