

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM112719\
 Data File : BM023954.D
 Acq On : 28 Nov 2019 16:04
 Operator : JU
 Sample : K5915-14
 Misc :
 ALS Vial : 38 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 JLNS2

Quant Time: Nov 28 21:37:26 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM112719MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Nov 27 17:16:22 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.49	152	459191	20.00	ng/ul	0.00
18) Naphthalene-d8	10.26	136	1836233	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.14	164	1098978	20.00	ng/ul	0.00
62) Phenanthrene-d10	16.90	188	2261704	20.00	ng/ul	0.00
78) Chrysene-d12	21.10	240	2229754	20.00	ng/ul	-0.01
86) Perylene-d12	23.25	264	2814565	20.00	ng/ul	-0.01

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.03	96	45955	4.45	ng/uL	0.00
5) Phenol-d5	6.67	99	857006	21.29	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.83	67	532121	22.73	ng/ul	0.00
9) 2-Chlorophenol-d4	7.02	132	703948	22.63	ng/ul	0.00
13) 4-Methylphenol-d8	8.20	113	675303	21.27	ng/ul	0.00
19) Nitrobenzene-d5	8.64	128	333116	23.37	ng/ul	0.00
22) 2-Nitrophenol-d4	9.36	143	363099	22.41	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	9.89	165	666492	21.73	ng/ul	0.00
29) 4-Chloroaniline-d4	10.40	131	787375	23.39	ng/ul	0.00
44) Dimethylphthalate-d6	13.57	166	2084593	24.41	ng/ul	0.00
47) Acenaphthylene-d8	13.83	160	2458303	24.56	ng/ul	0.00
52) 4-Nitrophenol-d4	14.37	143	251476	17.52	ng/ul	0.00
58) Fluorene-d10	15.14	176	1790824	24.92	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.29	200	243029	17.12	ng/ul	0.00
71) Anthracene-d10	17.00	188	2744783	25.41	ng/ul	0.00
79) Pyrene-d10	19.30	212	3035788	26.35	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.11	264	3885437	26.12	ng/ul	-0.01

Target Compounds					Ovalue	
6) Phenol	6.70	94	55875	1.350	ng/ul	93
45) Dimethylphthalate	13.62	163	913063	10.923	ng/ul	99

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

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