

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM112719\
 Data File : BM023929.D
 Acq On : 28 Nov 2019 01:09
 Operator : JU
 Sample : K5888-11
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 JLLR8

Quant Time: Nov 28 03:07:49 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM112719MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Nov 28 00:24:52 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.49	152	489491	20.00	ng/ul	0.00
18) Naphthalene-d8	10.26	136	2115426	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.15	164	1392800	20.00	ng/ul	0.00
62) Phenanthrene-d10	16.90	188	2932736	20.00	ng/ul	0.00
78) Chrysene-d12	21.11	240	2031869	20.00	ng/ul	0.00
86) Perylene-d12	23.25	264	2143317	20.00	ng/ul	-0.01

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.04	96	39478	3.59	ng/uL	0.00
5) Phenol-d5	6.68	99	832631	19.41	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.84	67	513150	20.57	ng/ul	0.00
9) 2-Chlorophenol-d4	7.03	132	663310	20.00	ng/ul	0.00
13) 4-Methylphenol-d8	8.20	113	672311	19.86	ng/ul	0.00
19) Nitrobenzene-d5	8.64	128	327286	19.93	ng/ul	0.00
22) 2-Nitrophenol-d4	9.36	143	368935	19.77	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	9.89	165	662959	18.76	ng/ul	0.00
29) 4-Chloroaniline-d4	10.41	131	845466	21.80	ng/ul	0.00
44) Dimethylphthalate-d6	13.57	166	2349957	21.71	ng/ul	0.00
47) Acenaphthylene-d8	13.84	160	2573331	20.29	ng/ul	0.00
52) 4-Nitrophenol-d4	14.38	143	279842	15.38	ng/ul	0.00
58) Fluorene-d10	15.15	176	1915008	21.03	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.29	200	292125	15.87	ng/ul	0.00
71) Anthracene-d10	17.00	188	3037341	21.68	ng/ul	0.00
79) Pyrene-d10	19.31	212	3100203	29.53	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.11	264	2584406	22.81	ng/ul	-0.01

Target Compounds				Ovalue		
6) Phenol	6.70	94	47046	1.066	ng/ul#	90
45) Dimethylphthalate	13.62	163	948091	8.950	ng/ul	100

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

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