

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM120319\
 Data File : BM023970.D
 Acq On : 03 Dec 2019 13:29
 Operator : JU
 Sample : K5914-02
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 JLNR3

Quant Time: Dec 03 14:54:42 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	444396	20.00	ng/ul	0.00
18) Naphthalene-d8	10.26	136	1930812	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.14	164	1208164	20.00	ng/ul	0.00
62) Phenanthrene-d10	16.90	188	2276612	20.00	ng/ul	0.00
78) Chrysene-d12	21.10	240	1482993	20.00	ng/ul	0.00
86) Perylene-d12	23.25	264	1868701	20.00	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.04	96	40456	4.05	ng/uL	0.00
5) Phenol-d5	6.67	99	821933	21.10	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.83	67	508513	22.45	ng/ul	0.00
9) 2-Chlorophenol-d4	7.02	132	659466	21.91	ng/ul	0.00
13) 4-Methylphenol-d8	8.20	113	667551	21.72	ng/ul	0.00
19) Nitrobenzene-d5	8.64	128	333466	22.25	ng/ul	0.00
22) 2-Nitrophenol-d4	9.35	143	367634	21.58	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	9.89	165	677513	21.00	ng/ul	0.00
29) 4-Chloroaniline-d4	10.40	131	782133	22.09	ng/ul	0.00
44) Dimethylphthalate-d6	13.57	166	2181862	23.24	ng/ul	0.00
47) Acenaphthylene-d8	13.83	160	2594659	23.58	ng/ul	0.00
52) 4-Nitrophenol-d4	14.37	143	257118	16.29	ng/ul	0.00
58) Fluorene-d10	15.14	176	1853892	23.47	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.29	200	235213	16.46	ng/ul	0.00
71) Anthracene-d10	17.00	188	2618096	24.08	ng/ul	0.00
79) Pyrene-d10	19.30	212	2347921	30.64	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.11	264	2547592	25.79	ng/ul	0.00
Target Compounds						
6) Phenol	6.70	94	74470	1.859	ng/ul	98
45) Dimethylphthalate	13.62	163	914999	9.957	ng/ul	100

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

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