

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM100319\
 Data File : BM022955.D
 Acq On : 04 Oct 2019 13:41
 Operator : JU
 Sample : K5058-06
 Misc :
 ALS Vial : 41 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DBAL9

Quant Time: Oct 04 17:05:36 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	223435	20.00	ng/ul	0.00
18) Naphthalene-d8	10.58	136	969531	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.43	164	598795	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.17	188	1132305	20.00	ng/ul	0.00
78) Chrysene-d12	21.34	240	807550	20.00	ng/ul	0.00
86) Perylene-d12	23.61	264	893567	20.00	ng/ul	0.00

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.26	96	24669	4.77	ng/uL	0.00
5) Phenol-d5	6.96	99	74522	3.78	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.13	67	142965	13.23	ng/ul	0.00
9) 2-Chlorophenol-d4	7.33	132	185547	11.67	ng/ul	0.00
13) 4-Methylphenol-d8	8.50	113	129195	7.99	ng/ul	0.00
19) Nitrobenzene-d5	8.95	128	102271	13.67	ng/ul	0.00
22) 2-Nitrophenol-d4	9.66	143	108470	13.28	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.20	165	215562	13.25	ng/ul	0.00
29) 4-Chloroaniline-d4	10.72	131	21941	1.11	ng/ul	0.00
44) Dimethylphthalate-d6	13.84	166	672573	14.47	ng/ul	0.00
47) Acenaphthylene-d8	14.12	160	765371	14.26	ng/ul	0.00
52) 4-Nitrophenol-d4	14.63	143	22115	2.41	ng/ul	0.00
58) Fluorene-d10	15.42	176	562830	14.21	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.53	200	73543	9.98	ng/ul	0.00
71) Anthracene-d10	17.26	188	828870	14.99	ng/ul	0.00
79) Pyrene-d10	19.56	212	787871	18.13	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.46	264	686828	14.86	ng/ul	0.00

Target Compounds					Ovalue	
6) Phenol	6.99	94	55085	2.662	ng/ul	97
45) Dimethylphthalate	13.88	163	262472	5.498	ng/ul	99

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

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