

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM100119\
 Data File : BM022895.D
 Acq On : 02 Oct 2019 16:31
 Operator : JU
 Sample : PB123400BL
 Misc :
 ALS Vial : 41 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :

Quant Time: Oct 03 00:59:01 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM092719MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Sep 27 18:03:27 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.80	152	227271	20.00	ng/ul	-0.01
18) Naphthalene-d8	10.59	136	1019079	20.00	ng/ul	-0.01
36) Acenaphthene-d10	14.43	164	623285	20.00	ng/ul	-0.02
62) Phenanthrene-d10	17.17	188	1099818	20.00	ng/ul	-0.01
78) Chrysene-d12	21.35	240	815628	20.00	ng/ul	-0.02
86) Perylene-d12	23.61	264	931107	20.00	ng/ul	-0.02

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	3.27	96	34704	6.60	ng/uL	0.00
5) Phenol-d5	6.97	99	611484	30.45	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.13	67	358272	32.60	ng/ul	-0.01
9) 2-Chlorophenol-d4	7.33	132	504036	31.17	ng/ul	-0.01
13) 4-Methylphenol-d8	8.50	113	508538	30.93	ng/ul	-0.01
19) Nitrobenzene-d5	8.95	128	252624	32.12	ng/ul	-0.01
22) 2-Nitrophenol-d4	9.67	143	287843	33.52	ng/ul	-0.01
26) 2,4-Dichlorophenol-d3	10.21	165	524577	30.68	ng/ul	-0.02
29) 4-Chloroaniline-d4	10.72	131	744948	35.86	ng/ul	-0.01
44) Dimethylphthalate-d6	13.84	166	1571072	32.48	ng/ul	-0.01
47) Acenaphthylene-d8	14.13	160	1819236	32.57	ng/ul	-0.01
52) 4-Nitrophenol-d4	14.63	143	218115	22.82	ng/ul	-0.01
58) Fluorene-d10	15.43	176	1269638	30.80	ng/ul	-0.01
63) 4,6-Dinitro-2-methylphenol	15.54	200	134485	18.79	ng/ul	-0.02
71) Anthracene-d10	17.27	188	1750188	32.59	ng/ul	-0.01
79) Pyrene-d10	19.56	212	1626381	37.05	ng/ul	-0.02
90) Benzo(a)pyrene-d12	23.47	264	1578040	32.76	ng/ul	-0.02

Target Compounds	Ovalue

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

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