

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM122319\
 Data File : BM024218.D
 Acq On : 23 Dec 2019 16:35
 Operator : JU
 Sample : PB125403BL
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK03

Quant Time: Dec 23 17:25:41 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM121719MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Dec 21 00:34:28 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.43	152	230227	20.00	ng/ul	0.00
18) Naphthalene-d8	10.20	136	1015832	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.10	164	642556	20.00	ng/ul	0.00
62) Phenanthrene-d10	16.85	188	1350422	20.00	ng/ul	0.00
78) Chrysene-d12	21.07	240	1245907	20.00	ng/ul	0.00
86) Perylene-d12	23.20	264	1470953	20.00	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	2.99	96	43384	8.32	ng/uL	0.00
5) Phenol-d5	6.63	99	707776	32.47	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.79	67	516509	39.51	ng/ul	0.00
9) 2-Chlorophenol-d4	6.97	132	564653	36.32	ng/ul	0.00
13) 4-Methylphenol-d8	8.16	113	580804	32.59	ng/ul	0.00
19) Nitrobenzene-d5	8.59	128	278327	38.13	ng/ul	0.00
22) 2-Nitrophenol-d4	9.30	143	303985	37.88	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	9.84	165	537071	34.42	ng/ul	0.00
29) 4-Chloroaniline-d4	10.36	131	769083	40.58	ng/ul	0.00
44) Dimethylphthalate-d6	13.52	166	1917881	40.36	ng/ul	0.00
47) Acenaphthylene-d8	13.79	160	2172332	38.15	ng/ul	0.00
52) 4-Nitrophenol-d4	14.34	143	225669	26.45	ng/ul	0.00
58) Fluorene-d10	15.10	176	1626427	39.05	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.25	200	223797	27.17	ng/ul	0.00
71) Anthracene-d10	16.95	188	2459930	40.04	ng/ul	0.00
79) Pyrene-d10	19.26	212	2603157	43.86	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.07	264	2919260	39.95	ng/ul	0.00

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

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