

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM090819\
 Data File : BM022599.D
 Acq On : 09 Sep 2019 11:48
 Operator : HP/JU
 Sample : PB122879BL
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK79

Quant Time: Sep 09 14:02:56 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM090519MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Sep 09 07:45:00 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.85	152	408466	20.00	ng/ul	0.00
18) Naphthalene-d8	10.65	136	1829082	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.49	164	1204600	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.22	188	2488092	20.00	ng/ul	0.00
78) Chrysene-d12	21.40	240	2559034	20.00	ng/ul	0.00
86) Perylene-d12	23.69	264	2799473	20.00	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	3.30	96	58155	5.93	ng/uL	0.00
5) Phenol-d5	7.01	99	1083232	28.75	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.18	67	656298	30.39	ng/ul	0.00
9) 2-Chlorophenol-d4	7.39	132	892084	29.97	ng/ul	0.00
13) 4-Methylphenol-d8	8.55	113	897205	29.29	ng/ul	0.00
19) Nitrobenzene-d5	9.00	128	443290	33.59	ng/ul	0.00
22) 2-Nitrophenol-d4	9.73	143	459855	37.80	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.26	165	888816	28.77	ng/ul	0.00
29) 4-Chloroaniline-d4	10.77	131	1346379	32.99	ng/ul	0.00
44) Dimethylphthalate-d6	13.89	166	2981941	30.55	ng/ul	0.00
47) Acenaphthylene-d8	14.18	160	3425330	29.31	ng/ul	0.00
52) 4-Nitrophenol-d4	14.66	143	539463	32.36	ng/ul	0.00
58) Fluorene-d10	15.47	176	2493994	30.52	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.58	200	330069	29.19	ng/ul	0.00
71) Anthracene-d10	17.32	188	3948894	30.98	ng/ul	0.00
79) Pyrene-d10	19.61	212	4453323	31.36	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.54	264	4811878	31.98	ng/ul	0.00

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

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