

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM091118\
 Data File : BM016723.D
 Acq On : 11 Sep 2018 15:29
 Operator : SJ/JU
 Sample : MDL-S-BL
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 MDL-S-BL

Quant Time: Sep 11 16:14:17 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM091018MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Sep 10 15:28:07 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.74	152	236882	20.00	ng/ul	0.00
18) Naphthalene-d8	10.52	136	992941	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.37	164	535336	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.12	188	1158834	20.00	ng/ul	0.00
78) Chrysene-d12	21.31	240	1233843	20.00	ng/ul	-0.01
86) Perylene-d12	23.56	264	1293386	20.00	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	3.28	96	33596	6.06	ng/uL	0.00
5) Phenol-d5	6.90	99	575760	28.82	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.08	67	372807	30.13	ng/ul	0.00
9) 2-Chlorophenol-d4	7.27	132	484636	29.72	ng/ul	0.00
13) 4-Methylphenol-d8	8.43	113	453765	29.35	ng/ul	-0.01
19) Nitrobenzene-d5	8.89	128	210710	32.29	ng/ul	0.00
22) 2-Nitrophenol-d4	9.60	143	194758	32.84	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.14	165	424403	28.16	ng/ul	0.00
29) 4-Chloroaniline-d4	10.66	131	645925	35.50	ng/ul	0.00
44) Dimethylphthalate-d6	13.79	166	1296364	30.16	ng/ul	0.00
47) Acenaphthylene-d8	14.07	160	1644017	29.61	ng/ul	0.00
52) 4-Nitrophenol-d4	14.57	143	170170	24.12	ng/ul	0.00
58) Fluorene-d10	15.37	176	1141323	30.42	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.49	200	98869	20.78	ng/ul	0.00
71) Anthracene-d10	17.22	188	1697496	30.36	ng/ul	0.00
79) Pyrene-d10	19.52	212	1839603	30.73	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.41	264	2057605	30.50	ng/ul	0.00

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

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