

Data Path : Z:\HPCHEM1\BNA_M\Data\BM091117\
 Data File : BM011497.D
 Acq On : 11 Sep 2017 15:26
 Operator : SJ/JU
 Sample : MDL-S-ME-BL
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 MDL-S-ME-BL

Quant Time: Sep 11 16:19:20 2017
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM090817MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Sep 11 14:52:48 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.98	152	486802	20.00	ng/ul	0.00
18) Naphthalene-d8	10.79	136	2184838	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.61	164	1294946	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.34	188	2888472	20.00	ng/ul	0.00
78) Chrysene-d12	21.51	240	3209677	20.00	ng/ul	0.00
86) Perylene-d12	23.89	264	3075232	20.00	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	3.39	96	97808	9.04	ng/uL	0.00
5) Phenol-d5	7.13	99	1851408	39.61	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.30	67	1379248	43.78	ng/ul	0.00
9) 2-Chlorophenol-d4	7.50	132	1504965	42.72	ng/ul	0.00
13) 4-Methylphenol-d8	8.67	113	1472188	40.34	ng/ul	0.00
19) Nitrobenzene-d5	9.14	128	745536	44.95	ng/ul	0.00
22) 2-Nitrophenol-d4	9.86	143	774636	44.94	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.40	165	1388125	40.00	ng/ul	0.00
29) 4-Chloroaniline-d4	10.92	131	2157175	56.32	ng/ul	0.00
44) Dimethylphthalate-d6	14.01	166	5088692	46.48	ng/ul	0.00
47) Acenaphthylene-d8	14.30	160	5852481	44.20	ng/ul	0.00
52) 4-Nitrophenol-d4	14.79	143	785740	38.19	ng/ul	0.00
58) Fluorene-d10	15.60	176	4282224	45.16	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.70	200	635204	37.79	ng/ul	0.00
71) Anthracene-d10	17.44	188	6943171	48.81	ng/ul	0.00
79) Pyrene-d10	19.73	212	7723256	51.39	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.74	264	7262361	49.55	ng/ul	0.00

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

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