

Data Path : Z:\HPCHEM1\BNA M\DATA\BM121115\
 Data File : BM003487.D
 Acq On : 11 Dec 2015 18:18
 Operator : UM/IZ
 Sample : G4718-04
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0AD2

Quant Time: Dec 12 00:57:34 2015
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-BM120115.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Dec 12 00:54:45 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.72	152	81807	20.00	ng/ul	0.00
18) Naphthalene-d8	10.51	136	359556	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.37	164	188155	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.12	188	376713	20.00	ng/ul	0.00
78) Chrysene-d12	21.32	240	382011	20.00	ng/ul	0.00
86) Perylene-d12	23.57	264	383459	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.18	96	2023	1.23	ng/uL	0.00
5) Phenol-d5	6.89	99	37894	5.80	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.05	67	90267	25.05	ng/ul	0.00
9) 2-Chlorophenol-d4	7.25	132	114176	21.14	ng/ul	0.00
13) 4-Methylphenol-d8	8.42	113	70664	12.90	ng/ul	0.00
19) Nitrobenzene-d5	8.87	128	64150	26.16	ng/ul	0.00
22) 2-Nitrophenol-d4	9.60	143	67211	26.21	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.13	165	115201	22.66	ng/ul	0.00
29) 4-Chloroaniline-d4	10.65	131	1467	0.26	ng/ul	0.00
44) Dimethylphthalate-d6	13.78	166	373433	26.40	ng/ul	0.00
47) Acenaphthylene-d8	14.06	160	463141	26.64	ng/ul	0.00
52) 4-Nitrophenol-d4	14.58	143	8731	3.44	ng/ul	0.00
58) Fluorene-d10	15.37	176	332923	27.35	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.49	200	36259	18.81	ng/ul	0.00
71) Anthracene-d10	17.22	188	518087	30.96	ng/ul	0.00
79) Pyrene-d10	19.52	212	547823	29.29	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.43	264	547609	29.60	ng/ul	0.00

Target Compounds Ovalue

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

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