

Data Path : Z:\HPCHEM1\BNA M\DATA\BM122115\
 Data File : BM003696.D
 Acq On : 22 Dec 2015 03:05
 Operator : UM/SJ
 Sample : G4867-19DL 50X
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 G9DB4DL

Quant Time: Dec 22 04:57:36 2015
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-BM121915.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Dec 22 02:40:18 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.69	152	26862	20.00	ng/ul	0.00
18) Naphthalene-d8	10.48	136	109485	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.35	164	58599	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.09	188	129682	20.00	ng/ul	0.00
78) Chrysene-d12	21.29	240	155007	20.00	ng/ul	0.00
86) Perylene-d12	23.53	264	158831	20.00	ng/ul	-0.01

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	0.00	96	0	0.00	ng/uL	
5) Phenol-d5	0.00	99	0	0.00	ng/ul	
7) Bis-(2-Chloroethyl)ether-d	7.03	67	309	0.26	ng/ul	0.00
9) 2-Chlorophenol-d4	7.23	132	551	0.30	ng/ul	0.00
13) 4-Methylphenol-d8	0.00	113	0	0.00	ng/ul	
19) Nitrobenzene-d5	0.00	128	0	0.00	ng/ul	
22) 2-Nitrophenol-d4	0.00	143	0	0.00	ng/ul	
26) 2,4-Dichlorophenol-d3	0.00	165	0	0.00	ng/ul	
29) 4-Chloroaniline-d4	0.00	131	0	0.00	ng/ul	
44) Dimethylphthalate-d6	13.76	166	2477	0.51	ng/ul	0.00
47) Acenaphthylene-d8	14.04	160	3052	0.55	ng/ul	0.00
52) 4-Nitrophenol-d4	0.00	143	0	0.00	ng/ul	
58) Fluorene-d10	15.35	176	2777	0.69	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	0.00	200	0	0.00	ng/ul	
71) Anthracene-d10	17.19	188	4136	0.70	ng/ul	0.00
79) Pyrene-d10	19.49	212	4211	0.62	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.39	264	4823	0.61	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Ovalue
28) Naphthalene	10.53	128	183074	32.46	ng/ul	100
34) 2-Methylnaphthalene	12.15	142	8437	2.00	ng/ul	95

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

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