

Data Path : Z:\HPCHEM1\BNA M\DATA\BM122115\
 Data File : BM003689.D
 Acq On : 21 Dec 2015 22:54
 Operator : UM/SJ
 Sample : G4867-14
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 G9DA9

Quant Time: Dec 22 03:03:44 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	28582	20.00	ng/ul	0.00
18) Naphthalene-d8	10.48	136	116965	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.34	164	62366	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.09	188	135759	20.00	ng/ul	0.00
78) Chrysene-d12	21.30	240	154337	20.00	ng/ul	0.00
86) Perylene-d12	23.54	264	159228	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.15	96	713	1.34	ng/uL	0.00
5) Phenol-d5	6.87	99	13089	5.43	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.03	67	38333	29.87	ng/ul	0.00
9) 2-Chlorophenol-d4	7.23	132	47614	24.22	ng/ul	0.00
13) 4-Methylphenol-d8	8.40	113	26052	12.49	ng/ul	0.00
19) Nitrobenzene-d5	8.85	128	27684	32.63	ng/ul	0.00
22) 2-Nitrophenol-d4	9.57	143	30318	31.76	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.11	165	48541	27.86	ng/ul	0.00
29) 4-Chloroaniline-d4	0.00	131	0d	0.00	ng/ul	
44) Dimethylphthalate-d6	13.76	166	166357	32.27	ng/ul	0.00
47) Acenaphthylene-d8	14.04	160	202225	34.46	ng/ul	0.00
52) 4-Nitrophenol-d4	14.56	143	3014	3.00	ng/ul	0.00
58) Fluorene-d10	15.34	176	144519	33.89	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.47	200	20737	27.10	ng/ul	0.00
71) Anthracene-d10	17.19	188	218905	35.44	ng/ul	0.00
79) Pyrene-d10	19.50	212	245814	36.46	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.40	264	276293	34.93	ng/ul	0.00

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

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