

Data Path : Z:\HPCHEM1\BNA M\DATA\BM040716\
 Data File : BM004919.D
 Acq On : 07 Apr 2016 16:56
 Operator : UM/SJ
 Sample : H2275-02
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 D9T75

Quant Time: Apr 08 01:10:32 2016
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-BM040116.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Apr 08 00:57:37 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.82	152	37122	20.00	ng/ul	0.00
18) Naphthalene-d8	10.61	136	175578	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.46	164	111764	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.20	188	281031	20.00	ng/ul	0.00
78) Chrysene-d12	21.39	240	375723	20.00	ng/ul	0.00
86) Perylene-d12	23.69	264	382269	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.32	96	744	0.81	ng/uL	0.00
5) Phenol-d5	6.99	99	15961	4.84	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.15	67	49274	27.57	ng/ul	0.00
9) 2-Chlorophenol-d4	7.35	132	52839	21.11	ng/ul	0.00
13) 4-Methylphenol-d8	8.52	113	33386	12.05	ng/ul	0.00
19) Nitrobenzene-d5	8.97	128	34033	27.32	ng/ul	0.00
22) 2-Nitrophenol-d4	9.70	143	37741	27.25	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.23	165	62692	24.28	ng/ul	0.00
29) 4-Chloroaniline-d4	10.76	131	264	0.09	ng/ul	0.02
44) Dimethylphthalate-d6	13.86	166	274816	31.17	ng/ul	0.00
47) Acenaphthylene-d8	14.15	160	312410	29.62	ng/ul	0.00
52) 4-Nitrophenol-d4	14.65	143	6932	3.98	ng/ul	0.00
58) Fluorene-d10	15.45	176	254905	33.45	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.57	200	45788	27.19	ng/ul	0.00
71) Anthracene-d10	17.30	188	455385	36.21	ng/ul	0.00
79) Pyrene-d10	19.60	212	554128	33.97	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.54	264	614528	36.28	ng/ul	0.00

Target Compounds Ovalue

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

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