

Data Path : Z:\HPCHEM1\BNA F\DATA\BF040216\
 Data File : BF085985.D
 Acq On : 2 Apr 2016 2:35
 Operator : UM/SJ
 Sample : H1824-05
 Misc : LOD 2PPM
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 LOD-WATER2016-01

Manual Integrations
 APPROVED

Sohil
 4/4/2016 7:43:30 PM

Quant Time: Apr 02 03:14:42 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF040216.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Apr 01 23:17:06 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.77	152	106794	20.00	ng	0.00
21) Naphthalene-d8	8.05	136	429469	20.00	ng	-0.01
38) Acenaphthene-d10	9.80	164	195276	20.00	ng	-0.01
63) Phenanthrene-d10	11.29	188	398217	20.00	ng	-0.01
75) Chrysene-d12	13.93	240	361335	20.00	ng	0.00
86) Perylene-d12	15.33	264	313488	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.37	112	767346	122.82	ng	0.00
7) Phenol-d6	6.42	99	1029218	129.69	ng	0.00
23) Nitrobenzene-d5	7.34	82	623970	85.68	ng	0.00
41) 2,4,6-Tribromophenol	10.60	330	286943	156.56	ng	0.00
44) 2-Fluorobiphenyl	9.14	172	1093003	93.06	ng	0.00
78) Terphenyl-d14	12.88	244	1085183	70.60	ng	0.00
Target Compounds						
3) Pyridine	3.46	79	6003m	0.91	ng	Qvalue 92
4) n-Nitrosodimethylamine	3.24	42	1696m	0.58	ng	
55) 4-Nitrophenol	10.01	139	2257m	1.04	ng	
64) 4,6-Dinitro-2-methylphenol	10.43	198	884	0.37	ng	
69) Pentachlorophenol	11.12	266	884	0.45	ng	

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

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