

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF061915\
 Data File : BF080061.D
 Acq On : 19 Jun 2015 17:00
 Operator : TP/IZ
 Sample : G2653-05
 Misc : LOD-SVOC-WATER-5PPM
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 LOD-WATER2015-01

Manual Integrations
 APPROVED

apatel
 6/22/2015 2:40:24 PM

Quant Time: Jun 22 12:51:05 2015
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF061715.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jun 17 16:41:53 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.33	152	39902	20.00	ng	-0.01
21) Naphthalene-d8	8.62	136	168655	20.00	ng	-0.01
38) Acenaphthene-d10	10.39	164	92091	20.00	ng	-0.01
63) Phenanthrene-d10	11.90	188	194580	20.00	ng	-0.01
75) Chrysene-d12	14.87	240	209372	20.00	ng	0.01
86) Perylene-d12	16.75	264	200494	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.94	112	265732	117.89	ng	0.00
7) Phenol-d6	6.93	99	329599	109.88	ng	-0.01
23) Nitrobenzene-d5	7.89	82	201299	69.48	ng	-0.01
41) 2,4,6-Tribromophenol	11.19	330	93993	104.38	ng	-0.01
44) 2-Fluorobiphenyl	9.69	172	420758	65.16	ng	-0.01
78) Terphenyl-d14	13.59	244	592707	66.11	ng	0.00
Target Compounds						
2) 1,4-Dioxane	3.25	88	2641	2.46	ng	# 73
3) Pyridine	4.15	79	6733	2.36	ng	# 80
4) n-Nitrosodimethylamine	3.99	42	4306	3.18	ng	# 64
25) Isophorone	8.14	82	17772	3.05	ng	# 84
32) Benzoic acid	8.28	122	755m	0.48	ng	
40) Hexachlorocyclopentadiene	9.50	237	2317	7.80	ng	82
53) 2,4-Dinitrophenol	10.44	184	362	6.43	ng	# 1
64) 4,6-Dinitro-2-methylphenol	10.97	198	1373	5.25	ng	# 77

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

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