

Data Path : Z:\HPCHEM1\BNA G\DATA\BG022416\
 Data File : BG021059.D
 Acq On : 24 Feb 2016 19:52
 Operator : UM/SJ
 Sample : H1520-19
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 40050+15(0-2)

Manual Integrations
 APPROVED

Sohil
 2/25/2016 2:40:15 PM

Quant Time: Feb 25 00:31:27 2016
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG022316.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Feb 24 00:35:06 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.21	152	4131	20.00	ng	0.00
21) Naphthalene-d8	11.03	136	22542	20.00	ng	0.00
38) Acenaphthene-d10	14.82	164	12376	20.00	ng	0.00
63) Phenanthrene-d10	17.55	188	23526	20.00	ng	0.00
75) Chrysene-d12	21.82	240	17537	20.00	ng	0.00
86) Perylene-d12	25.14	264	17182	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.78	112	39632	168.20	ng	-0.01
7) Phenol-d6	7.41	99	61856	170.46	ng	0.00
23) Nitrobenzene-d5	9.39	82	39837	103.76	ng	0.00
41) 2,4,6-Tribromophenol	16.31	330	22064	148.82	ng	0.00
44) 2-Fluorobiphenyl	13.45	172	83247	93.64	ng	0.00
78) Terphenyl-d14	20.14	244	70612	52.25	ng	0.00
Target Compounds						
20) 3+4-Methylphenols	9.01	107	1258	3.52	ng	94
49) Dimethylphthalate	14.27	163	17274	17.12	ng	98
51) Acenaphthene	14.88	154	2320	2.94	ng	96
70) Phenanthrene	17.60	178	25293	19.01	ng	99
71) Anthracene	17.69	178	6007	4.53	ng	98
74) Fluoranthene	19.59	202	29826	22.29	ng	98
77) Pyrene	19.95	202	22690	11.32	ng	98
80) Benzo(a)anthracene	21.80	228	9816	9.33	ng	96
82) Chrysene	21.87	228	8429	8.59	ng	95
87) Benzo(b)fluoranthene	24.09	252	11199m	9.89	ng	
88) Benzo(k)fluoranthene	24.15	252	3133	2.93	ng	# 92
89) Benzo(a)pyrene	24.98	252	7929	7.81	ng	# 94

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

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