

Data Path : S:\HPCHEM1\BNA_G\DATA\BG072015\
 Data File : BG017988.D
 Acq On : 20 Jul 2015 14:40
 Operator : TP/UM
 Sample : G2976-07
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 BRIDGE14-8(0-2)

Manual Integrations
 APPROVED

apatel
 7/22/2015 9:43:26 AM

Quant Time: Jul 20 18:36:53 2015
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG071715.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Jul 18 01:50:30 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.59	152	156395	20.00	ng	0.00
21) Naphthalene-d8	10.37	136	667012	20.00	ng	0.00
38) Acenaphthene-d10	14.25	164	393913	20.00	ng	0.00
63) Phenanthrene-d10	17.00	188	741652	20.00	ng	0.00
75) Chrysene-d12	21.20	240	766068	20.00	ng	0.00
86) Perylene-d12	23.42	264	781969	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.20	112	934071	104.19	ng	0.00
7) Phenol-d6	6.78	99	1212965	103.27	ng	0.01
23) Nitrobenzene-d5	8.74	82	721164	70.03	ng	0.00
41) 2,4,6-Tribromophenol	15.74	330	431946	108.23	ng	0.00
44) 2-Fluorobiphenyl	12.86	172	1493266	65.69	ng	0.00
78) Terphenyl-d14	19.65	244	1747742	57.47	ng	0.00
Target Compounds						Qvalue
49) Dimethylphthalate	13.71	163	525873	19.22	ng	99
70) Phenanthrene	17.04	178	379915	9.93	ng	97
71) Anthracene	17.13	178	45052	1.21	ng	99
74) Fluoranthene	19.06	202	585378	13.98	ng	97
77) Pyrene	19.43	202	524181	12.27	ng	97
80) Benzo(a)anthracene	21.19	228	222405	5.50	ng	99
82) Chrysene	21.23	228	251350	6.52	ng	95
85) Indeno(1,2,3-cd)pyrene	25.66	276	127563	2.74	ng	# 89
87) Benzo(b)fluoranthene	22.75	252	317368m	7.42	ng	
88) Benzo(k)fluoranthene	22.79	252	88083m	2.06	ng	
89) Benzo(a)pyrene	23.32	252	214599	5.38	ng	99
91) Benzo(g,h,i)perylene	26.36	276	120253	3.02	ng	99

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

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