

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG060815\
 Data File : BG017548.D
 Acq On : 8 Jun 2015 18:04
 Operator : TP/IZ
 Sample : G2555-05 5X
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 DUSO-005

Manual Integrations
 APPROVED

apatel
 6/9/2015 3:56:33 PM

Quant Time: Jun 09 03:29:52 2015
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG060315.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 09 02:10:45 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.61	152	19237	20.00	ng	0.00
21) Naphthalene-d8	10.40	136	78623	20.00	ng	0.00
38) Acenaphthene-d10	14.28	164	54726	20.00	ng	0.00
63) Phenanthrene-d10	17.04	188	115307	20.00	ng	0.00
75) Chrysene-d12	21.23	240	130809	20.00	ng	0.00
86) Perylene-d12	23.46	264	133238	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.20	112	29286	26.82	ng	0.00
7) Phenol-d6	6.82	99	35780	24.05	ng	0.00
23) Nitrobenzene-d5	8.78	82	27255	17.30	ng	0.00
41) 2,4,6-Tribromophenol	15.79	330	17041	22.42	ng	0.00
44) 2-Fluorobiphenyl	12.89	172	58156	16.03	ng	0.00
78) Terphenyl-d14	19.67	244	85140	13.45	ng	0.00
Target Compounds						Qvalue
57) Fluorene	15.34	166	10691	2.30	ng	97
70) Phenanthrene	17.08	178	92651	14.41	ng	97
71) Anthracene	17.17	178	14123m	2.20	ng	
74) Fluoranthene	19.10	202	111104	13.22	ng	97
77) Pyrene	19.47	202	146409	18.17	ng	98
80) Benzo(a)anthracene	21.21	228	67934	8.44	ng	96
82) Chrysene	21.27	228	67898	9.32	ng	96
85) Indeno(1,2,3-cd)pyrene	25.74	276	27535	3.01	ng	# 90
87) Benzo(b)fluoranthene	22.80	252	64545m	7.64	ng	
88) Benzo(k)fluoranthene	22.84	252	21709m	2.69	ng	
89) Benzo(a)pyrene	23.37	252	57251	7.39	ng	96
91) Benzo(g,h,i)perylene	26.44	276	23496	3.07	ng	97

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

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