

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE070815\
 Data File : BE090073.D
 Acq On : 8 Jul 2015 23:55
 Operator : TP/IZ
 Sample : G2782-01 5X
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 X3SS0330006-150624

Manual Integrations
 APPROVED

apatel
 7/9/2015 2:21:12 PM

Quant Time: Jul 09 02:25:06 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\8270-SIM-BE062915.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 30 07:01:01 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.62	152	18904	5.00	ng	0.00
6) Naphthalene-d8	10.37	136	85461	5.00	ng	-0.02
12) Acenaphthene-d10	14.23	164	46121	5.00	ng	0.00
18) Phenanthrene-d10	16.96	188	101374	5.00	ng	-0.02
25) Chrysene-d12	21.12	240	118911	5.00	ng	-0.01
32) Perylene-d12	23.44	264	113522	5.00	ng	-0.02
System Monitoring Compounds						
4) 2-Fluorophenol	5.25	112	3410	0.79	ng	0.00
5) Phenol-d6	6.80	99	5424	0.89	ng	0.00
7) Nitrobenzene-d5	8.76	82	3133	0.46	ng	0.00
13) 2,4,6-Tribromophenol	15.73	330	1183	0.88	ng	0.00
14) 2-Fluorobiphenyl	12.85	172	7011	0.51	ng	0.00
27) Terphenyl-d14	19.58	244	9847	0.50	ng	-0.01
Target Compounds						Qvalue
15) Acenaphthylene	13.94	152	4012	0.05	ng	99
22) Phenanthrene	17.00	178	1812m	0.07	ng	
23) Anthracene	17.10	178	1223	0.05	ng	99
24) Fluoranthene	19.00	202	6073	0.21	ng	100
26) Pyrene	19.37	202	5872	0.21	ng	# 96
28) Benzo(a)anthracene	21.10	228	4754	0.16	ng	91
29) Chrysene	21.16	228	5392m	0.17	ng	
31) Indeno(1,2,3-cd)pyrene	25.83	276	4204	0.14	ng	99
33) Benzo(b)fluoranthene	22.73	252	8198	0.33	ng	# 85
34) Benzo(k)fluoranthene	22.77	252	2694m	0.10	ng	
35) Benzo(a)pyrene	23.33	252	5125	0.22	ng	# 82
37) Benzo(g,h,i)perylene	26.57	276	4816	0.20	ng	# 88

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

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