

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE070815\
 Data File : BE090067.D
 Acq On : 8 Jul 2015 20:20
 Operator : TP/IZ
 Sample : G2824-13
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 X1SS0820006-150626

Manual Integrations
 APPROVED

apatel
 7/9/2015 2:20:56 PM

Quant Time: Jul 09 01:45:23 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	7.61	152	21467	5.00	ng	0.00
6) Naphthalene-d8	10.37	136	94134	5.00	ng	-0.01
12) Acenaphthene-d10	14.22	164	52183	5.00	ng	-0.01
18) Phenanthrene-d10	16.96	188	116827	5.00	ng	-0.02
25) Chrysene-d12	21.12	240	135877	5.00	ng	-0.01
32) Perylene-d12	23.43	264	124894	5.00	ng	-0.02
System Monitoring Compounds						
4) 2-Fluorophenol	5.25	112	23998	4.87	ng	0.00
5) Phenol-d6	6.80	99	41109	5.96	ng	0.00
7) Nitrobenzene-d5	8.75	82	23995	3.21	ng	0.00
13) 2,4,6-Tribromophenol	15.73	330	10087	6.60	ng	0.00
14) 2-Fluorobiphenyl	12.85	172	58358	3.72	ng	0.00
27) Terphenyl-d14	19.58	244	95182	4.20	ng	-0.02
Target Compounds						Qvalue
22) Phenanthrene	17.00	178	6317	0.21	ng	97
23) Anthracene	17.08	178	902m	0.03	ng	
24) Fluoranthene	19.00	202	22040	0.66	ng	99
26) Pyrene	19.36	202	20833	0.64	ng	97
28) Benzo(a)anthracene	21.10	228	13626	0.40	ng	98
29) Chrysene	21.15	228	18050	0.51	ng	97
31) Indeno(1,2,3-cd)pyrene	25.83	276	10406	0.31	ng	96
33) Benzo(b)fluoranthene	22.73	252	22982	0.83	ng	# 96
34) Benzo(k)fluoranthene	22.77	252	9029m	0.29	ng	
35) Benzo(a)pyrene	23.32	252	14420	0.57	ng	96
36) Dibenzo(a,h)anthracene	25.85	278	2485	0.10	ng	# 78
37) Benzo(g,h,i)perylene	26.57	276	11251	0.42	ng	94

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

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