

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE063015\
 Data File : BE090033.D
 Acq On : 1 Jul 2015 5:59
 Operator : TP/IZ
 Sample : G2811-01DL 20X
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 X1SS0520006-150625-DDL

Quant Time: Jul 01 08:46:14 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	16340	5.00	ng	0.00
6) Naphthalene-d8	10.37	136	70464	5.00	ng	-0.01
12) Acenaphthene-d10	14.23	164	38800	5.00	ng	0.00
18) Phenanthrene-d10	16.96	188	86290	5.00	ng	-0.02
25) Chrysene-d12	21.12	240	109103	5.00	ng	-0.01
32) Perylene-d12	23.44	264	106742	5.00	ng	-0.02
System Monitoring Compounds						
4) 2-Fluorophenol	5.25	112	869	0.23	ng	0.00
5) Phenol-d6	6.80	99	1207	0.23	ng	0.00
7) Nitrobenzene-d5	8.76	82	766	0.14	ng	0.00
13) 2,4,6-Tribromophenol	15.73	330	212	0.19	ng	0.00
14) 2-Fluorobiphenyl	12.85	172	1826	0.16	ng	0.00
27) Terphenyl-d14	19.58	244	2712	0.15	ng	0.00
Target Compounds						Qvalue
9) Naphthalene	10.42	128	1030	0.06	ng	# 83
16) Acenaphthene	14.29	154	901	0.09	ng	98
22) Phenanthrene	17.00	178	28905	1.31	ng	100
23) Anthracene	17.10	178	5726m	0.28	ng	
24) Fluoranthene	19.01	202	111630	4.52	ng	100
26) Pyrene	19.37	202	105988	4.08	ng	# 96
28) Benzo(a)anthracene	21.11	228	82930	3.04	ng	99
29) Chrysene	21.15	228	89800	3.17	ng	98
31) Indeno(1,2,3-cd)pyrene	25.84	276	49661	1.83	ng	97
33) Benzo(b)fluoranthene	22.73	252	117897	5.01	ng	99
34) Benzo(k)fluoranthene	22.77	252	44688m	1.70	ng	
35) Benzo(a)pyrene	23.33	252	80473	3.70	ng	98
36) Dibenzo(a,h)anthracene	25.85	278	12385	0.55	ng	95
37) Benzo(g,h,i)perylene	26.58	276	51797	2.29	ng	98

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

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