

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE063015\
 Data File : BE090032.D
 Acq On : 1 Jul 2015 5:20
 Operator : TP/IZ
 Sample : G2824-02 2X
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 X1SS0650006-150626-D

Quant Time: Jul 01 17:16:18 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	14907	5.00	ng	0.00
6) Naphthalene-d8	10.37	136	64220	5.00	ng	-0.01
12) Acenaphthene-d10	14.23	164	34621	5.00	ng	0.00
18) Phenanthrene-d10	16.96	188	80425	5.00	ng	-0.02
25) Chrysene-d12	21.12	240	103552	5.00	ng	-0.01
32) Perylene-d12	23.44	264	103327	5.00	ng	-0.02

System Monitoring Compounds						
4) 2-Fluorophenol	5.24	112	12336	3.60	ng	0.00
5) Phenol-d6	6.80	99	18257	3.81	ng	0.00
7) Nitrobenzene-d5	8.76	82	10547	2.07	ng	0.00
13) 2,4,6-Tribromophenol	15.73	330	3560	3.51	ng	0.00
14) 2-Fluorobiphenyl	12.85	172	23201	2.23	ng	-0.01
27) Terphenyl-d14	19.58	244	36626	2.12	ng	-0.01

Target Compounds						Qvalue
9) Naphthalene	10.42	128	2094	0.14	ng	# 93
11) 2-Methylnaphthalene	12.04	142	342	0.04	ng	# 87
16) Acenaphthene	14.28	154	1426	0.16	ng	93
17) Fluorene	15.27	166	886	0.07	ng	# 93
22) Phenanthrene	17.01	178	52474	2.55	ng	100
23) Anthracene	17.10	178	10116m	0.54	ng	
24) Fluoranthene	19.01	202	168817	7.34	ng	99
26) Pyrene	19.37	202	159253	6.45	ng	96
28) Benzo(a)anthracene	21.10	228	122798	4.75	ng	99
29) Chrysene	21.16	228	133163	4.96	ng	98
31) Indeno(1,2,3-cd)pyrene	25.84	276	83880	3.26	ng	96
33) Benzo(b)fluoranthene	22.74	252	182019	7.99	ng	98
34) Benzo(k)fluoranthene	22.77	252	73086m	2.87	ng	
35) Benzo(a)pyrene	23.33	252	124544	5.92	ng	98
36) Dibenzo(a,h)anthracene	25.85	278	20425	0.95	ng	97
37) Benzo(g,h,i)perylene	26.58	276	86034	3.93	ng	98

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

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