

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE070215\
 Data File : BE090053.D
 Acq On : 2 Jul 2015 16:33
 Operator : TP/IZ
 Sample : G2810-02RE
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 X1SS0600624-150625RE

Manual Integrations
 APPROVED

apatel
 7/6/2015 2:41:17 PM

Quant Time: Jul 03 01:45:05 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	11663	5.00	ng	-0.02
6) Naphthalene-d8	10.37	136	53719	5.00	ng	-0.02
12) Acenaphthene-d10	14.22	164	30576	5.00	ng	-0.01
18) Phenanthrene-d10	16.96	188	68800	5.00	ng	-0.02
25) Chrysene-d12	21.12	240	81273	5.00	ng	-0.01
32) Perylene-d12	23.44	264	74836	5.00	ng	-0.02
System Monitoring Compounds						
4) 2-Fluorophenol	5.24	112	19859	7.41	ng	-0.01
5) Phenol-d6	6.79	99	31652	8.44	ng	-0.01
7) Nitrobenzene-d5	8.75	82	18677	4.38	ng	-0.01
13) 2,4,6-Tribromophenol	15.73	330	5321	5.94	ng	0.00
14) 2-Fluorobiphenyl	12.84	172	41363	4.50	ng	-0.01
27) Terphenyl-d14	19.58	244	59168	4.37	ng	-0.01
Target Compounds						Qvalue
9) Naphthalene	10.41	128	323	0.03	ng	# 57
11) 2-Methylnaphthalene	12.03	142	265	0.03	ng	# 87
15) Acenaphthylene	13.94	152	2187	0.04	ng	97
16) Acenaphthene	14.28	154	387	0.05	ng	91
17) Fluorene	15.27	166	822	0.08	ng	# 85
21) Pentachlorophenol	16.63	266	305	0.45	ng	92
22) Phenanthrene	17.00	178	5538	0.31	ng	98
23) Anthracene	17.10	178	1605m	0.10	ng	
24) Fluoranthene	19.00	202	12232	0.62	ng	100
26) Pyrene	19.36	202	11644	0.60	ng	97
28) Benzo(a)anthracene	21.10	228	8302	0.41	ng	99
29) Chrysene	21.16	228	9108	0.43	ng	98
30) Bis(2-ethylhexyl)phthalate	21.05	149	1013	0.41	ng	100
31) Indeno(1,2,3-cd)pyrene	25.84	276	4334	0.21	ng	99
33) Benzo(b)fluoranthene	22.73	252	10083	0.61	ng	96
34) Benzo(k)fluoranthene	22.77	252	4222m	0.23	ng	
35) Benzo(a)pyrene	23.33	252	6740	0.44	ng	# 94
36) Dibenzo(a,h)anthracene	25.85	278	1608	0.10	ng	# 79
37) Benzo(g,h,i)perylene	26.57	276	4305	0.27	ng	94

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

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