

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE070115\
 Data File : BE090043.D
 Acq On : 1 Jul 2015 15:05
 Operator : TP/IZ
 Sample : G2810-26DL 10X
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 X1SS0520006-150625DL

Manual Integrations
 APPROVED

apatel
 7/2/2015 3:35:45 PM

Quant Time: Jul 02 05:51:28 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	17764	5.00	ng	-0.01
6) Naphthalene-d8	10.37	136	78518	5.00	ng	-0.02
12) Acenaphthene-d10	14.22	164	45826	5.00	ng	-0.01
18) Phenanthrene-d10	16.96	188	106575	5.00	ng	-0.02
25) Chrysene-d12	21.12	240	132714	5.00	ng	-0.01
32) Perylene-d12	23.44	264	125802	5.00	ng	-0.02
System Monitoring Compounds						
4) 2-Fluorophenol	5.24	112	2212	0.54	ng	-0.01
5) Phenol-d6	6.79	99	3161	0.55	ng	-0.01
7) Nitrobenzene-d5	8.75	82	1947	0.31	ng	-0.01
13) 2,4,6-Tribromophenol	15.73	330	496	0.37	ng	0.00
14) 2-Fluorobiphenyl	12.85	172	4691	0.34	ng	0.00
27) Terphenyl-d14	19.58	244	7263	0.33	ng	0.00
Target Compounds						Qvalue
9) Naphthalene	10.41	128	1489	0.08	ng	# 90
16) Acenaphthene	14.29	154	984	0.08	ng	93
17) Fluorene	15.27	166	703	0.04	ng	# 81
22) Phenanthrene	17.00	178	25632	0.94	ng	99
23) Anthracene	17.10	178	4806m	0.19	ng	
24) Fluoranthene	19.00	202	87279	2.86	ng	100
26) Pyrene	19.36	202	82661	2.61	ng	97
28) Benzo(a)anthracene	21.11	228	62348	1.88	ng	97
29) Chrysene	21.15	228	68966	2.00	ng	98
31) Indeno(1,2,3-cd)pyrene	25.83	276	34751	1.05	ng	97
33) Benzo(b)fluoranthene	22.73	252	88158	3.18	ng	98
34) Benzo(k)fluoranthene	22.77	252	30892m	1.00	ng	
35) Benzo(a)pyrene	23.33	252	57556	2.25	ng	98
36) Dibenzo(a,h)anthracene	25.85	278	8846	0.34	ng	97
37) Benzo(g,h,i)perylene	26.57	276	35937	1.35	ng	98

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

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