

Data Path : Z:\HPCHEM1\BNA E\DATA\BE041116\
 Data File : BE091642.D
 Acq On : 11 Apr 2016 23:37
 Operator : UM/SJ
 Sample : H2090-05
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 KY038PS027-160330

Manual Integrations
 APPROVED

Sohil
 4/12/2016 5:33:53 PM

Quant Time: Apr 12 06:59:52 2016
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE040816.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Apr 08 16:48:14 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.80	152	196038	5.00	ng	-0.02
7) Naphthalene-d8	10.57	136	997052	5.00	ng	-0.02
13) Acenaphthene-d10	14.42	164	614477	5.00	ng	0.00
19) Phenanthrene-d10	17.16	188	1540667	5.00	ng	0.00
26) Chrysene-d12	21.31	240	1666874	5.00	ng	0.00
35) Perylene-d12	23.76	264	1564470	5.00	ng	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Fluorophenol	5.41	112	102720	2.20	ng	0.00
5) Phenol-d6	6.97	99	101369	1.63	ng	0.00
8) Nitrobenzene-d5	8.96	82	242954	4.19	ng	0.00
14) 2,4,6-Tribromophenol	15.90	330	195134	8.66	ng	-0.01
15) 2-Fluorobiphenyl	13.05	172	642204	3.76	ng	0.00
29) Terphenyl-d14	19.76	244	1119531	5.38	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
24) Anthracene	17.29	178	8152	0.03	ng	99
30) Benzo(a)anthracene	21.29	228	15322m	0.04	ng	
32) Chrysene	21.34	228	10380m	0.03	ng	
33) Bis(2-ethylhexyl)phthalate	21.22	149	35928	0.39	ng	# 83
34) Indeno(1,2,3-cd)pyrene	26.32	276	9909	0.03	ng	97
36) Benzo(b)fluoranthene	23.01	252	10102	0.03	ng	# 77
37) Benzo(k)fluoranthene	23.06	252	9635	0.03	ng	# 84
39) Dibenzo(a,h)anthracene	26.34	278	9358	0.03	ng	94
40) Benzo(g,h,i)perylene	27.10	276	8931	0.03	ng	96

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

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