

Data Path : Z:\HPCHEM1\BNA E\DATA\BE041116\
 Data File : BE091635.D
 Acq On : 11 Apr 2016 19:17
 Operator : UM/SJ
 Sample : PB89409BSD
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 PB89409BSD

Quant Time: Apr 12 01:48:00 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	0.00	152	0	0.00	ng	-7.82
7) Naphthalene-d8	10.63	136	9	5.00	ng	0.04
13) Acenaphthene-d10	14.51	164	2	5.00	ng	0.09
19) Phenanthrene-d10	17.33	188	3	5.00	ng	0.17
26) Chrysene-d12	0.00	240	0	0.00	ng	-21.31
35) Perylene-d12	23.77	264	327	5.00	ng	0.01
System Monitoring Compounds						
4) 2-Fluorophenol	5.44	112	17	0.00	ng	0.03
5) Phenol-d6	7.04	99	7	0.00	ng	0.07
8) Nitrobenzene-d5	8.98	82	4	7.65	ng	0.02
14) 2,4,6-Tribromophenol	0.00	330	0	0.00	ng	
15) 2-Fluorobiphenyl	13.10	172	5	8.99	ng	0.06
29) Terphenyl-d14	19.78	244	105	0.00	ng	0.02
Target Compounds						
9) Nitrobenzene	9.06	77	8	11.92	ng	# 22
10) Naphthalene	10.63	128	16	8.62	ng	# 1
12) 2-Methylnaphthalene	12.29	142	6	5.06	ng	# 48
16) Acenaphthylene	14.27	152	5	6.44	ng	# 14
17) Acenaphthene	14.57	154	5	10.19	ng	# 1
18) Fluorene	15.60	166	7	11.41	ng	# 7
20) 4-Bromophenyl-phenylether	16.52	248	1	9.45	ng	# 1
21) Hexachlorobenzene	16.66	284	22	200.36	ng	# 50
23) Phenanthrene	17.33	178	28	40.89	ng	# 54
24) Anthracene	17.44	178	35	57.68	ng	# 1
25) Fluoranthene	19.40	202	16	22.08	ng	# 63
36) Benzo(b)fluoranthene	23.02	252	301	3.96	ng	# 1
37) Benzo(k)fluoranthene	23.07	252	359	4.79	ng	# 1
38) Benzo(a)pyrene	23.67	252	304	4.31	ng	# 1
39) Dibenzo(a,h)anthracene	26.38	278	694	9.99	ng	# 39
40) Benzo(g,h,i)perylene	27.12	276	384	5.37	ng	# 28

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

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