

Data Path : Z:\HPCHEM1\BNA F\DATA\BF110716\
 Data File : BF091064.D
 Acq On : 7 Nov 2016 19:01
 Operator : UM/SJ
 Sample : H5543-03
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-110-0.5-1.0

Manual Integrations
 APPROVED

sohil
 11/8/2016 4:32:10 PM

Quant Time: Nov 08 11:39:49 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF110316.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 08 00:06:22 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.67	152	197922	20.00	ng	0.00
21) Naphthalene-d8	7.96	136	737577	20.00	ng	0.00
38) Acenaphthene-d10	9.71	164	316590	20.00	ng	0.00
63) Phenanthrene-d10	11.22	188	249035	20.00	ng	0.02
75) Chrysene-d12	13.92	240	118642	20.00	ng	0.09
86) Perylene-d12	15.32	264	246476	20.00	ng	0.11
System Monitoring Compounds						
5) 2-Fluorophenol	5.25	112	1218708	101.69	ng	0.00
7) Phenol-d6	6.33	99	1614027	99.23	ng	0.00
23) Nitrobenzene-d5	7.24	82	1029965	76.18	ng	0.00
41) 2,4,6-Tribromophenol	10.51	330	202967	70.91	ng	0.00
44) 2-Fluorobiphenyl	9.04	172	1444988	78.58	ng	-0.01
78) Terphenyl-d14	12.84	244	295125	62.07	ng	0.06
Target Compounds						
31) Naphthalene	7.97	128	73344	2.08	ng	98
48) Acenaphthylene	9.57	152	67853	2.48	ng	# 70
49) Dimethylphthalate	9.45	163	524090	25.25	ng	# 98
70) Phenanthrene	11.24	178	162808	12.30	ng	# 66
74) Fluoranthene	12.47	202	97453	7.10	ng	70
77) Pyrene	12.69	202	100971	13.00	ng	# 74
85) Indeno(1,2,3-cd)pyrene	16.63	276	117619	21.34	ng	# 89
91) Benzo(q,h,i)perylene	17.00	276	109757	9.65	ng	94

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

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