

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF091818\
 Data File : BF109106.D
 Acq On : 18 Sep 2018 20:02
 Operator : JU/SJ
 Sample : J4952-01
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 2015-161-20-RB36584RT-RB47184R

Manual Integrations
 APPROVED

Sohil
 9/19/2018 4:23:47 PM

Quant Time: Sep 19 05:07:39 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF091418.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Sep 14 13:26:52 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.72	152	128833	20.00	ng	0.00
21) Naphthalene-d8	8.00	136	466878	20.00	ng	0.00
38) Acenaphthene-d10	9.75	164	200347	20.00	ng	0.00
63) Phenanthrene-d10	11.23	188	340110	20.00	ng	0.00
75) Chrysene-d12	13.85	240	352465	20.00	ng	0.00
86) Perylene-d12	15.26	264	348539	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.33	112	524046	67.56	ng	0.00
7) Phenol-d6	6.35	99	677511	67.74	ng	0.00
23) Nitrobenzene-d5	7.28	82	418157	50.24	ng	-0.01
41) 2,4,6-Tribromophenol	10.54	330	146332	67.30	ng	0.00
44) 2-Fluorobiphenyl	9.08	172	759843	59.42	ng	0.00
78) Terphenyl-d14	12.81	244	689332	46.36	ng	0.00
Target Compounds						
10) Phenol	6.37	94	41540	3.691	ng	93
49) Dimethylphthalate	9.47	163	119094	8.317	ng	98
74) Fluoranthene	12.44	202	38874	2.209	ng	95
83) Bis(2-ethylhexyl)phthalate	13.85	149	75310	5.319	ng	98
85) Indeno(1,2,3-cd)pyrene	16.61	276	50607m	2.965	ng	
87) Benzo(b)fluoranthene	14.86	252	96783m	5.019	ng	
89) Benzo(a)pyrene	15.20	252	71995	4.204	ng	# 96
91) Benzo(g,h,i)perylene	17.01	276	49257	3.425	ng	93

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

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