

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF091417\
 Data File : BF098547.D
 Acq On : 15 Sep 2017 00:14
 Operator : SJ/JU
 Sample : I5161-17
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 RL-7-0-11

Manual Integrations
 APPROVED

Sohil
 9/15/2017 5:44:26 PM

Quant Time: Sep 15 07:00:29 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF091117.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Sep 12 02:14:50 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.65	152	149185	20.00	ng	-0.02
21) Naphthalene-d8	7.93	136	605943	20.00	ng	-0.02
38) Acenaphthene-d10	9.68	164	259860	20.00	ng	-0.02
63) Phenanthrene-d10	11.16	188	396605	20.00	ng	-0.02
75) Chrysene-d12	13.80	240	295854	20.00	ng	-0.02
86) Perylene-d12	15.19	264	309136	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.26	112	1189671	117.90	ng	-0.01
7) Phenol-d6	6.30	99	1520557	118.69	ng	-0.02
23) Nitrobenzene-d5	7.22	82	926152	85.21	ng	-0.02
41) 2,4,6-Tribromophenol	10.47	330	220900	127.37	ng	-0.02
44) 2-Fluorobiphenyl	9.01	172	1489607	97.16	ng	-0.02
78) Terphenyl-d14	12.75	244	953654	69.58	ng	-0.02
Target Compounds						
10) Phenol	6.32	94	38898	2.614	ng	86
49) Dimethylphthalate	9.40	163	213156	10.597	ng	100
70) Phenanthrene	11.19	178	49577	2.358	ng	97
74) Fluoranthene	12.37	202	83347	3.960	ng	99
77) Pyrene	12.60	202	75114	3.218	ng	98
80) Benzo(a)anthracene	13.79	228	35165	2.027	ng	# 89
82) Chrysene	13.82	228	41263	2.353	ng	95
87) Benzo(b)fluoranthene	14.80	252	56326m	2.898	ng	
89) Benzo(a)pyrene	15.14	252	40256	2.325	ng	# 89

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

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