

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF101818\
 Data File : BF109980.D
 Acq On : 19 Oct 2018 1:02
 Operator : JU/SJ
 Sample : J5570-02
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TS-07

Manual Integrations
 APPROVED

Sohil
 10/19/2018 12:15:03 PM

Quant Time: Oct 19 05:36:44 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF101518.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Oct 16 14:58:24 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.97	152	229103	20.00	ng	0.00
21) Naphthalene-d8	8.26	136	872265	20.00	ng	-0.01
39) Acenaphthene-d10	10.02	164	386096	20.00	ng	-0.01
64) Phenanthrene-d10	11.51	188	617949	20.00	ng	0.00
76) Chrysene-d12	14.15	240	432967	20.00	ng	-0.01
87) Perylene-d12	15.68	264	412376	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.61	112	1738374	120.28	ng	0.01
7) Phenol-d6	6.60	99	2199802	122.57	ng	0.00
23) Nitrobenzene-d5	7.54	82	1419518	109.63	ng	-0.01
42) 2,4,6-Tribromophenol	10.81	330	455613	136.30	ng	0.00
45) 2-Fluorobiphenyl	9.33	172	2107873	87.11	ng	0.00
79) Terphenyl-d14	13.10	244	1506687	73.08	ng	0.00
Target Compounds						
10) Phenol	6.62	94	245562	12.673	ng	Qvalue # 59
50) Dimethylphthalate	9.72	163	362916	13.275	ng	100
71) Phenanthrene	11.53	178	77592	2.425	ng	99
75) Fluoranthene	12.72	202	168489	5.272	ng	99
78) Pyrene	12.95	202	133825	4.208	ng	98
81) Benzo(a)anthracene	14.14	228	52125	2.044	ng	99
83) Chrysene	14.17	228	67407	2.777	ng	98
88) Benzo(b)fluoranthene	15.22	252	88485m	3.722	ng	
90) Benzo(a)pyrene	15.61	252	54162	2.477	ng	# 97

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

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