

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF101818\
 Data File : BF109981.D
 Acq On : 19 Oct 2018 1:29
 Operator : JU/SJ
 Sample : J5570-03
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TS-08

Manual Integrations
 APPROVED

Sohil
 10/19/2018 12:15:06 PM

Quant Time: Oct 19 11:23:06 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	210129	20.00	ng	0.00
21) Naphthalene-d8	8.26	136	821464	20.00	ng	-0.01
39) Acenaphthene-d10	10.02	164	367237	20.00	ng	-0.01
64) Phenanthrene-d10	11.51	188	587064	20.00	ng	0.00
76) Chrysene-d12	14.15	240	411906	20.00	ng	-0.01
87) Perylene-d12	15.68	264	416459	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.60	112	1086302	81.95	ng	0.00
7) Phenol-d6	6.60	99	1353976	82.25	ng	0.00
23) Nitrobenzene-d5	7.54	82	846202	69.39	ng	-0.01
42) 2,4,6-Tribromophenol	10.81	330	275354	86.60	ng	0.00
45) 2-Fluorobiphenyl	9.33	172	1332718	57.91	ng	0.00
79) Terphenyl-d14	13.09	244	946628	48.26	ng	-0.01
Target Compounds						
10) Phenol	6.61	94	150462	8.466	ng	Qvalue # 59
50) Dimethylphthalate	9.72	163	224064	8.617	ng	99
75) Fluoranthene	12.72	202	100811	3.320	ng	97
78) Pyrene	12.95	202	75975	2.511	ng	98
88) Benzo(b)fluoranthene	15.22	252	52957m	2.206	ng	

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

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