

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF110918\
 Data File : BF110616.D
 Acq On : 9 Nov 2018 10:04
 Operator : JU/SJ
 Sample : J5813-03
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TS-C

Manual Integrations
 APPROVED

Sohil
 11/12/2018 10:29:51 AM

Quant Time: Nov 09 15:50:44 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF110818.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Nov 09 15:13:13 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.95	152	78826	20.00	ng	0.00
21) Naphthalene-d8	8.23	136	314260	20.00	ng	0.00
39) Acenaphthene-d10	9.99	164	138911	20.00	ng	0.00
64) Phenanthrene-d10	11.47	188	216414	20.00	ng	-0.01
76) Chrysene-d12	14.13	240	155265	20.00	ng	-0.01
87) Perylene-d12	15.65	264	141210	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.57	112	500321	103.10	ng	0.00
7) Phenol-d6	6.57	99	638534	102.90	ng	0.00
23) Nitrobenzene-d5	7.51	82	402820	73.82	ng	0.00
42) 2,4,6-Tribromophenol	10.77	330	89342	63.83	ng	-0.01
45) 2-Fluorobiphenyl	9.30	172	656071	67.45	ng	0.00
79) Terphenyl-d14	13.06	244	409582	49.40	ng	0.00
Target Compounds						
10) Phenol	6.58	94	41717	5.797	ng	Qvalue # 69
50) Dimethylphthalate	9.69	163	106208	10.247	ng	98
75) Fluoranthene	12.70	202	31336	2.696	ng	98
78) Pyrene	12.93	202	27043	2.262	ng	99
88) Benzo(b)fluoranthene	15.20	252	19525m	2.311	ng	

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

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