

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF110918\
 Data File : BF110641.D
 Acq On : 9 Nov 2018 21:36
 Operator : JU/SJ
 Sample : J5813-06
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TS-F

Manual Integrations
 APPROVED

Sohil
 11/12/2018 10:30:13 AM

Quant Time: Nov 11 03:49:56 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	85074	20.00	ng	0.00
21) Naphthalene-d8	8.23	136	363706	20.00	ng	0.00
39) Acenaphthene-d10	9.99	164	173613	20.00	ng	0.00
64) Phenanthrene-d10	11.48	188	309582	20.00	ng	0.00
76) Chrysene-d12	14.13	240	176575	20.00	ng	-0.01
87) Perylene-d12	15.66	264	162832	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.56	112	466698	89.11	ng	0.00
7) Phenol-d6	6.57	99	611395	91.29	ng	0.00
23) Nitrobenzene-d5	7.51	82	385516	61.04	ng	0.00
42) 2,4,6-Tribromophenol	10.78	330	142653	81.55	ng	0.00
45) 2-Fluorobiphenyl	9.30	172	707442	58.20	ng	0.00
79) Terphenyl-d14	13.06	244	539009	57.17	ng	0.00
Target Compounds						
10) Phenol	6.58	94	41578	5.354	ng	Qvalue # 64
50) Dimethylphthalate	9.69	163	104336	8.054	ng	99
71) Phenanthrene	11.50	178	43008	2.593	ng	97
75) Fluoranthene	12.70	202	81513	4.902	ng	98
78) Pyrene	12.93	202	62273	4.580	ng	97
83) Chrysene	14.16	228	25985	2.584	ng	98
88) Benzo(b)fluoranthene	15.20	252	34518m	3.544	ng	
90) Benzo(a)pyrene	15.59	252	19963	2.256	ng	# 95

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

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