

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF110918\
 Data File : BF110643.D
 Acq On : 9 Nov 2018 22:31
 Operator : JU/SJ
 Sample : J5813-09
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TS-I

Manual Integrations
 APPROVED

Sohil
 11/12/2018 10:30:19 AM

Quant Time: Nov 11 03:53:56 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	101246	20.00	ng	0.00
21) Naphthalene-d8	8.23	136	403387	20.00	ng	0.00
39) Acenaphthene-d10	9.99	164	170371	20.00	ng	0.00
64) Phenanthrene-d10	11.48	188	260202	20.00	ng	0.00
76) Chrysene-d12	14.13	240	187419	20.00	ng	-0.01
87) Perylene-d12	15.66	264	139801	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.57	112	643584	103.25	ng	0.00
7) Phenol-d6	6.57	99	823785	103.35	ng	0.00
23) Nitrobenzene-d5	7.51	82	524246	74.84	ng	0.00
42) 2,4,6-Tribromophenol	10.78	330	149130	86.87	ng	0.00
45) 2-Fluorobiphenyl	9.30	172	853081	71.51	ng	0.00
79) Terphenyl-d14	13.06	244	555006	55.46	ng	0.00
Target Compounds						
10) Phenol	6.59	94	53784	5.819	ng	Qvalue # 64
50) Dimethylphthalate	9.69	163	126082	9.918	ng	99
71) Phenanthrene	11.50	178	68490	4.913	ng	98
75) Fluoranthene	12.70	202	145340	10.399	ng	98
78) Pyrene	12.93	202	112958	7.828	ng	98
81) Benzo(a)anthracene	14.12	228	48500	4.329	ng	97
83) Chrysene	14.16	228	60401	5.660	ng	96
86) Indeno(1,2,3-cd)pyrene	17.22	276	23125	3.020	ng	# 89
88) Benzo(b)fluoranthene	15.20	252	63710m	7.618	ng	
89) Benzo(k)fluoranthene	15.23	252	20389m	2.467	ng	
90) Benzo(a)pyrene	15.59	252	33013	4.345	ng	# 94
92) Benzo(a,h,i)perylene	17.71	276	21082	3.304	ng	# 88

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

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