

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF110918\
 Data File : BF110642.D
 Acq On : 9 Nov 2018 22:03
 Operator : JU/SJ
 Sample : J5813-07
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TS-G

Manual Integrations
 APPROVED

Sohil
 11/12/2018 10:30:16 AM

Quant Time: Nov 11 03:52:11 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	84219	20.00	ng	0.00
21) Naphthalene-d8	8.23	136	348012	20.00	ng	0.00
39) Acenaphthene-d10	9.99	164	153351	20.00	ng	0.00
64) Phenanthrene-d10	11.48	188	238809	20.00	ng	0.00
76) Chrysene-d12	14.13	240	163216	20.00	ng	-0.01
87) Perylene-d12	15.66	264	129982	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.56	112	476221	91.85	ng	0.00
7) Phenol-d6	6.57	99	627034	94.57	ng	0.00
23) Nitrobenzene-d5	7.51	82	397354	65.75	ng	0.00
42) 2,4,6-Tribromophenol	10.78	330	119400	77.27	ng	0.00
45) 2-Fluorobiphenyl	9.30	172	672300	62.61	ng	0.00
79) Terphenyl-d14	13.06	244	436884	50.13	ng	0.00
Target Compounds						
10) Phenol	6.58	94	42200	5.489	ng	Qvalue # 66
50) Dimethylphthalate	9.69	163	105024	9.178	ng	99
71) Phenanthrene	11.50	178	28167	2.202	ng	99
75) Fluoranthene	12.70	202	59589	4.646	ng	96
78) Pyrene	12.93	202	50192	3.994	ng	98
83) Chrysene	14.16	228	23910	2.573	ng	# 95
88) Benzo(b)fluoranthene	15.20	252	30198m	3.884	ng	
90) Benzo(a)pyrene	15.59	252	16740	2.369	ng	# 89

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

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