

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF110818\
 Data File : BF110604.D
 Acq On : 9 Nov 2018 00:02
 Operator : JU/SJ
 Sample : J5768-03 2X
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 OK-02-110718-B

Manual Integrations
 APPROVED

Sohil
 11/9/2018 4:26:15 PM

Quant Time: Nov 09 13:24:18 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF110818.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Nov 08 14:14:47 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.95	152	79503	20.00	ng	0.00
21) Naphthalene-d8	8.23	136	334804	20.00	ng	0.00
39) Acenaphthene-d10	9.98	164	147681	20.00	ng	-0.01
64) Phenanthrene-d10	11.48	188	232108	20.00	ng	0.00
76) Chrysene-d12	14.13	240	173314	20.00	ng	-0.01
87) Perylene-d12	15.65	264	147235	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.56	112	192942	39.42	ng	0.00
7) Phenol-d6	6.56	99	250667	40.05	ng	-0.01
23) Nitrobenzene-d5	7.50	82	154998	26.66	ng	-0.01
42) 2,4,6-Tribromophenol	10.77	330	43564	29.28	ng	-0.01
45) 2-Fluorobiphenyl	9.30	172	293906	28.42	ng	0.00
79) Terphenyl-d14	13.06	244	196470	21.23	ng	-0.01
Target Compounds						Qvalue
71) Phenanthrene	11.50	178	116614	9.378	ng	98
72) Anthracene	11.56	178	26815m	2.138	ng	
75) Fluoranthene	12.70	202	91156	7.312	ng	98
78) Pyrene	12.93	202	89043	6.673	ng	98
81) Benzo(a)anthracene	14.12	228	35371	3.414	ng	97
83) Chrysene	14.15	228	32032m	3.246	ng	
88) Benzo(b)fluoranthene	15.20	252	27162m	3.084	ng	
90) Benzo(a)pyrene	15.59	252	23300	2.911	ng	95

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

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