

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF080118\
 Data File : BF107888.D
 Acq On : 1 Aug 2018 14:53
 Operator : JU/SJ
 Sample : J4241-01
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 FS-RO-226

Manual Integrations
 APPROVED

Sohil
 8/2/2018 9:17:02 AM

Quant Time: Aug 01 16:19:22 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF073018.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jul 30 17:48:53 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.97	152	115956	20.00	ng	0.00
21) Naphthalene-d8	8.24	136	499411	20.00	ng	0.00
38) Acenaphthene-d10	10.00	164	212232	20.00	ng	0.00
63) Phenanthrene-d10	11.50	188	399932	20.00	ng	0.00
75) Chrysene-d12	14.15	240	389823	20.00	ng	0.00
86) Perylene-d12	15.69	264	448754	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.60	112	546150	80.05	ng	0.00
7) Phenol-d6	6.60	99	714302	79.44	ng	0.00
23) Nitrobenzene-d5	7.53	82	442659	51.06	ng	0.00
41) 2,4,6-Tribromophenol	10.80	330	191799	66.51	ng	0.00
44) 2-Fluorobiphenyl	9.31	172	763304	50.12	ng	0.00
78) Terphenyl-d14	13.08	244	707826	39.37	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.71	163	141848	8.900	ng	98
70) Phenanthrene	11.52	178	350312	18.447	ng	97
71) Anthracene	11.58	178	132205	6.036	ng	97
74) Fluoranthene	12.72	202	462625	20.737	ng	96
77) Pyrene	12.95	202	521313	18.572	ng	97
80) Benzo(a)anthracene	14.14	228	242842	10.997	ng	97
82) Chrysene	14.18	228	198209	8.322	ng	99
85) Indeno(1,2,3-cd)pyrene	17.28	276	98591m	5.442	ng	
87) Benzo(b)fluoranthene	15.23	252	207645m	9.382	ng	
88) Benzo(k)fluoranthene	15.25	252	109972m	4.047	ng	
89) Benzo(a)pyrene	15.62	252	197195	8.631	ng	97
91) Benzo(a,h,i)perylene	17.75	276	92968	4.821	ng	95

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

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