

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF110618\
 Data File : BF110536.D
 Acq On : 7 Nov 2018 00:49
 Operator : JU/SJ
 Sample : J5764-15 5X
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 JC-03-110218-H

Manual Integrations
 APPROVED

Sohil
 11/7/2018 2:54:51 PM

Quant Time: Nov 07 13:10:12 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF102318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Nov 01 16:06:12 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.96	152	77489	20.00	ng	0.00
21) Naphthalene-d8	8.24	136	286299	20.00	ng	0.00
39) Acenaphthene-d10	10.00	164	107434	20.00	ng	0.00
64) Phenanthrene-d10	11.49	188	189969	20.00	ng	0.00
76) Chrysene-d12	14.14	240	128763	20.00	ng	0.00
87) Perylene-d12	15.67	264	97019	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.57	112	77238	16.23	ng	0.00
7) Phenol-d6	6.57	99	100839	16.57	ng	0.00
23) Nitrobenzene-d5	7.52	82	53023	10.98	ng	0.00
42) 2,4,6-Tribromophenol	10.79	330	16003	15.04	ng	0.00
45) 2-Fluorobiphenyl	9.31	172	108299	15.83	ng	0.00
79) Terphenyl-d14	13.07	244	95550	15.43	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.70	163	16179	2.040	ng	Qvalue # 98
71) Phenanthrene	11.52	178	57352	5.770	ng	99
75) Fluoranthene	12.71	202	81847	8.113	ng	95
78) Pyrene	12.95	202	74141	8.032	ng	98
81) Benzo(a)anthracene	14.13	228	31962	4.186	ng	# 92
83) Chrysene	14.17	228	29092	4.011	ng	# 93
84) Bis(2-ethylhexyl)phthalate	14.09	149	25874	4.777	ng	# 90
88) Benzo(b)fluoranthene	15.22	252	29094m	5.193	ng	
90) Benzo(a)pyrene	15.60	252	20630	4.002	ng	# 75
92) Benzo(g,h,i)perylene	17.73	276	12284	2.802	ng	# 85

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

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