

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF103018\
 Data File : BF110323.D
 Acq On : 31 Oct 2018 00:21
 Operator : JU/SJ
 Sample : J5685-03DL 2X
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 DOCUMENTATION-27DL

Manual Integrations
 APPROVED

Sohil
 10/31/2018 11:15:11 AM

Quant Time: Oct 31 07:50:59 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF102318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 29 17:29:00 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.97	152	89431	20.00	ng	0.00
21) Naphthalene-d8	8.26	136	331728	20.00	ng	0.00
39) Acenaphthene-d10	10.02	164	130989	20.00	ng	0.00
64) Phenanthrene-d10	11.51	188	223566	20.00	ng	0.00
76) Chrysene-d12	14.16	240	185443	20.00	ng	0.00
87) Perylene-d12	15.70	264	122760	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.59	112	244285	44.48	ng	0.00
7) Phenol-d6	6.59	99	296523	42.21	ng	0.00
23) Nitrobenzene-d5	7.53	82	174747	31.24	ng	0.00
42) 2,4,6-Tribromophenol	10.81	330	50635	39.02	ng	0.00
45) 2-Fluorobiphenyl	9.33	172	316997	38.00	ng	0.00
79) Terphenyl-d14	13.09	244	271940	30.50	ng	0.00
Target Compounds						
10) Phenol	6.60	94	19808	2.638	ng	Qvalue # 78
50) Dimethylphthalate	9.72	163	48328	4.998	ng	98
71) Phenanthrene	11.53	178	201375	17.214	ng	99
72) Anthracene	11.59	178	77224	6.586	ng	99
75) Fluoranthene	12.73	202	521007	43.885	ng	97
78) Pyrene	12.96	202	518469	38.998	ng	99
81) Benzo(a)anthracene	14.15	228	277650	25.247	ng	97
83) Chrysene	14.19	228	227340	21.765	ng	97
86) Indeno(1,2,3-cd)pyrene	17.28	276	103408	11.934	ng	98
88) Benzo(b)fluoranthene	15.24	252	271715m	38.329	ng	
89) Benzo(k)fluoranthene	15.27	252	73142m	10.495	ng	
90) Benzo(a)pyrene	15.63	252	175679	26.932	ng	99
91) Dibenzo(a,h)anthracene	17.29	278	26353	4.682	ng	# 85
92) Benzo(g,h,i)perylene	17.76	276	100608	18.140	ng	99

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

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