

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF102918\
 Data File : BF110285.D
 Acq On : 30 Oct 2018 1:48
 Operator : JU/SJ
 Sample : J5685-07
 Misc :
 ALS Vial : 35 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 DOCUMENTATION-31

Manual Integrations
 APPROVED

Sohil
 10/30/2018 3:04:29 PM

Quant Time: Oct 30 09:09:29 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	93013	20.00	ng	0.00
21) Naphthalene-d8	8.25	136	347368	20.00	ng	0.00
39) Acenaphthene-d10	10.01	164	135763	20.00	ng	0.00
64) Phenanthrene-d10	11.50	188	217681	20.00	ng	0.00
76) Chrysene-d12	14.16	240	143907	20.00	ng	0.00
87) Perylene-d12	15.69	264	104249	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.59	112	496803	86.98	ng	0.00
7) Phenol-d6	6.59	99	612990	83.91	ng	0.00
23) Nitrobenzene-d5	7.53	82	364971	62.32	ng	0.00
42) 2,4,6-Tribromophenol	10.80	330	90227	67.09	ng	0.00
45) 2-Fluorobiphenyl	9.33	172	684935	79.22	ng	0.00
79) Terphenyl-d14	13.09	244	518103	74.88	ng	0.00
Target Compounds						
10) Phenol	6.60	94	44044	5.640	ng	Qvalue # 64
50) Dimethylphthalate	9.72	163	112419	11.218	ng	100
71) Phenanthrene	11.53	178	230474	20.235	ng	98
72) Anthracene	11.58	178	53494	4.686	ng	99
75) Fluoranthene	12.73	202	398940	34.512	ng	99
78) Pyrene	12.96	202	331078	32.091	ng	98
81) Benzo(a)anthracene	14.14	228	145813	17.086	ng	99
83) Chrysene	14.18	228	131618	16.238	ng	98
86) Indeno(1,2,3-cd)pyrene	17.26	276	56886	8.460	ng	98
88) Benzo(b)fluoranthene	15.23	252	140074m	23.268	ng	
89) Benzo(k)fluoranthene	15.26	252	47473m	8.021	ng	
90) Benzo(a)pyrene	15.62	252	91618	16.539	ng	# 96
91) Dibenzo(a,h)anthracene	17.26	278	15369	3.215	ng	# 79
92) Benzo(a,h,i)perylene	17.74	276	55876	11.863	ng	97

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

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