

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF102918\
 Data File : BF110287.D
 Acq On : 30 Oct 2018 2:42
 Operator : JU/SJ
 Sample : J5685-02
 Misc :
 ALS Vial : 37 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 DOCUMENTATION-26

Manual Integrations
 APPROVED

Sohil
 10/30/2018 3:04:34 PM

Quant Time: Oct 30 09:15:25 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	104979	20.00	ng	0.00
21) Naphthalene-d8	8.25	136	425009	20.00	ng	0.00
39) Acenaphthene-d10	10.02	164	187551	20.00	ng	0.00
64) Phenanthrene-d10	11.51	188	308283	20.00	ng	0.00
76) Chrysene-d12	14.16	240	198969	20.00	ng	0.00
87) Perylene-d12	15.69	264	161803	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.59	112	574847	89.17	ng	0.00
7) Phenol-d6	6.59	99	727041	88.17	ng	0.00
23) Nitrobenzene-d5	7.53	82	449157	62.68	ng	0.00
42) 2,4,6-Tribromophenol	10.81	330	139466	75.07	ng	0.00
45) 2-Fluorobiphenyl	9.33	172	807356	67.60	ng	0.00
79) Terphenyl-d14	13.09	244	596544	62.35	ng	0.00
Target Compounds						
10) Phenol	6.60	94	50254	5.702	ng	Qvalue # 70
31) Naphthalene	8.27	128	56922	2.906	ng	97
50) Dimethylphthalate	9.72	163	141973	10.255	ng	100
71) Phenanthrene	11.53	178	193700	12.008	ng	99
72) Anthracene	11.59	178	54095	3.346	ng	98
75) Fluoranthene	12.73	202	269514	16.463	ng	98
78) Pyrene	12.96	202	233762	16.388	ng	98
81) Benzo(a)anthracene	14.15	228	119870	10.159	ng	98
83) Chrysene	14.19	228	101320	9.041	ng	97
86) Indeno(1,2,3-cd)pyrene	17.27	276	42196	4.539	ng	97
88) Benzo(b)fluoranthene	15.23	252	130323m	13.948	ng	
89) Benzo(k)fluoranthene	15.26	252	32224m	3.508	ng	
90) Benzo(a)pyrene	15.63	252	83259	9.684	ng	# 97
92) Benzo(a,h,i)perylene	17.75	276	39850	5.451	ng	95

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

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