

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF111418\
 Data File : BF110769.D
 Acq On : 14 Nov 2018 17:28
 Operator : JU/SJ
 Sample : J5889-01
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WALL-A

Manual Integrations
 APPROVED

Sohil
 11/15/2018 10:59:29 AM

Quant Time: Nov 15 04:42:57 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF110818.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 13 12:43:10 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.95	152	75309	20.00	ng	0.00
21) Naphthalene-d8	8.22	136	310782	20.00	ng	0.00
39) Acenaphthene-d10	9.99	164	143650	20.00	ng	0.00
64) Phenanthrene-d10	11.48	188	240496	20.00	ng	0.00
76) Chrysene-d12	14.13	240	154414	20.00	ng	-0.01
87) Perylene-d12	15.66	264	140846	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.57	112	544768	117.50	ng	0.00
7) Phenol-d6	6.57	99	696074	117.41	ng	0.00
23) Nitrobenzene-d5	7.51	82	438167	81.19	ng	0.00
42) 2,4,6-Tribromophenol	10.78	330	154076	106.45	ng	0.00
45) 2-Fluorobiphenyl	9.30	172	833502	82.87	ng	0.00
79) Terphenyl-d14	13.06	244	619006	75.07	ng	0.00
Target Compounds						
10) Phenol	6.58	94	40757	5.929	ng	Qvalue # 77
50) Dimethylphthalate	9.69	163	94822	8.846	ng	100
71) Phenanthrene	11.50	178	135613	10.525	ng	99
72) Anthracene	11.56	178	32060	2.467	ng	97
75) Fluoranthene	12.70	202	214872	16.634	ng	98
78) Pyrene	12.93	202	186550	15.690	ng	98
81) Benzo(a)anthracene	14.12	228	86949	9.421	ng	98
83) Chrysene	14.16	228	78181	8.891	ng	98
86) Indeno(1,2,3-cd)pyrene	17.23	276	39851	6.316	ng	96
88) Benzo(b)fluoranthene	15.20	252	107349m	12.740	ng	
89) Benzo(k)fluoranthene	15.23	252	32280m	3.877	ng	
90) Benzo(a)pyrene	15.59	252	77630	10.140	ng	98
92) Benzo(a,h,i)perylene	17.71	276	34991m	5.444	ng	

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

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