

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121718\
 Data File : BF111552.D
 Acq On : 17 Dec 2018 21:21
 Operator : JU/SJ
 Sample : J6403-13
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-2-B

Manual Integrations
 APPROVED

Sohil
 12/18/2018 12:49:29 PM

Quant Time: Dec 18 06:34:48 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF121418.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 14 13:26:17 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.79	152	134014	20.00	ng	0.00
21) Naphthalene-d8	8.07	136	809196	20.00	ng	-0.01
39) Acenaphthene-d10	9.82	164	364049	20.00	ng	-0.01
64) Phenanthrene-d10	11.31	188	566223	20.00	ng	0.00
76) Chrysene-d12	13.94	240	374802	20.00	ng	0.00
87) Perylene-d12	15.39	264	390806	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.42	112	961330	119.37	ng	0.00
7) Phenol-d6	6.44	99	1064014	99.32	ng	0.00
23) Nitrobenzene-d5	7.35	82	702630	51.71	ng	0.00
42) 2,4,6-Tribromophenol	10.62	330	297382	91.19	ng	0.00
45) 2-Fluorobiphenyl	9.15	172	1582184	67.13	ng	-0.01
79) Terphenyl-d14	12.89	244	1083744	67.90	ng	0.00
Target Compounds						
10) Phenol	6.45	94	68671	5.755	ng	95
50) Dimethylphthalate	9.54	163	191700	7.312	ng	97
71) Phenanthrene	11.33	178	448914	15.386	ng	93
72) Anthracene	11.38	178	128457m	4.305	ng	
75) Fluoranthene	12.52	202	504087	17.660	ng	97
78) Pyrene	12.75	202	446733	16.906	ng	94
81) Benzo(a)anthracene	13.93	228	212169	10.411	ng	99
83) Chrysene	13.97	228	193382	9.008	ng	96
86) Indeno(1,2,3-cd)pyrene	16.80	276	89595	5.782	ng	92
88) Benzo(b)fluoranthene	14.97	252	242322m	10.635	ng	
89) Benzo(k)fluoranthene	14.99	252	87226m	4.007	ng	
90) Benzo(a)pyrene	15.32	252	180421	8.781	ng	97
92) Benzo(a,h,i)perylene	17.22	276	82517	5.187	ng	98

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

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