

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF112318\
 Data File : BF110949.D
 Acq On : 23 Nov 2018 17:35
 Operator : JU/SJ
 Sample : J5999-09
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSSI-1116-S-AB-4

Manual Integrations
 APPROVED

MOHAMMAD
 11/26/2018 3:12:10 PM

Quant Time: Nov 23 21:28:30 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	6.93	152	65623	20.00	ng	-0.01
21) Naphthalene-d8	8.22	136	252331	20.00	ng	0.00
39) Acenaphthene-d10	9.97	164	99025	20.00	ng	-0.02
64) Phenanthrene-d10	11.47	188	168159	20.00	ng	-0.01
76) Chrysene-d12	14.13	240	158340	20.00	ng	-0.01
87) Perylene-d12	15.66	264	106658	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.56	112	392487	97.15	ng	0.00
7) Phenol-d6	6.56	99	485743	94.02	ng	-0.01
23) Nitrobenzene-d5	7.50	82	307004	70.07	ng	-0.01
42) 2,4,6-Tribromophenol	10.77	330	86055	86.24	ng	-0.01
45) 2-Fluorobiphenyl	9.29	172	404303	58.31	ng	-0.01
79) Terphenyl-d14	13.05	244	355705	42.07	ng	-0.02
Target Compounds						
10) Phenol	6.57	94	24852	4.149	ng	Qvalue # 75
50) Dimethylphthalate	9.69	163	29012	3.926	ng	100
71) Phenanthrene	11.50	178	32126	3.566	ng	98
75) Fluoranthene	12.69	202	84735	9.381	ng	98
78) Pyrene	12.93	202	73749	6.049	ng	97
81) Benzo(a)anthracene	14.12	228	44039	4.653	ng	97
83) Chrysene	14.16	228	40890	4.535	ng	97
86) Indeno(1,2,3-cd)pyrene	17.25	276	16729	2.586	ng	96
88) Benzo(b)fluoranthene	15.20	252	45551m	7.139	ng	
89) Benzo(k)fluoranthene	15.23	252	13371m	2.121	ng	
90) Benzo(a)pyrene	15.60	252	27912	4.815	ng	# 93
92) Benzo(a,h,i)perylene	17.74	276	16088	3.305	ng	# 91

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

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