

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF120719\
 Data File : BF118134.D
 Acq On : 7 Dec 2019 17:35
 Operator : JU
 Sample : K6147-01
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 P001-WC001

Manual Integrations
 APPROVED

mohammad
 12/10/2019 11:47:40 AM

Quant Time: Dec 09 07:03:32 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF120619.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 06 17:34:43 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.86	152	124686	20.00	ng	0.00
21) Naphthalene-d8	8.15	136	497774	20.00	ng	0.00
39) Acenaphthene-d10	9.91	164	269145	20.00	ng	-0.01
64) Phenanthrene-d10	11.40	188	472723	20.00	ng	0.00
76) Chrysene-d12	14.06	240	271538	20.00	ng	-0.01
87) Perylene-d12	15.54	264	298540	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.48	112	677513	95.17	ng	0.00
7) Phenol-d6	6.49	99	885279	102.81	ng	0.00
23) Nitrobenzene-d5	7.43	82	529855	59.88	ng	-0.01
42) 2,4,6-Tribromophenol	10.70	330	299315	91.54	ng	-0.01
45) 2-Fluorobiphenyl	9.23	172	1108308	62.66	ng	0.00
79) Terphenyl-d14	13.00	244	1066188	67.52	ng	0.00
Target Compounds						
17) 2-Methylphenol	7.12	107	15819	2.478	ng	95
50) Dimethylphthalate	9.62	163	127375	6.407	ng	100
71) Phenanthrene	11.43	178	465105	18.508	ng	99
72) Anthracene	11.48	178	52795	2.106	ng	98
75) Fluoranthene	12.62	202	552648	21.510	ng	98
78) Pyrene	12.85	202	464102	21.689	ng	99
81) Benzo(a)anthracene	14.04	228	193216	10.291	ng	99
83) Chrysene	14.08	228	195145	10.880	ng	97
86) Indeno(1,2,3-cd)pyrene	17.04	276	106942	7.087	ng	95
88) Benzo(b)fluoranthene	15.11	252	272228m	13.788	ng	
89) Benzo(k)fluoranthene	15.13	252	72347m	3.814	ng	
90) Benzo(a)pyrene	15.48	252	182672	10.475	ng	99
92) Benzo(a,h,i)perylene	17.50	276	99604	6.930	ng	96

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

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