

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF033120\
 Data File : BF119663.D
 Acq On : 31 Mar 2020 20:22
 Operator : MA/SJ
 Sample : L2055-08
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-7-(1-3)

Manual Integrations
 APPROVED

mohammad
 4/1/2020 11:55:51 AM

Quant Time: Apr 01 00:56:41 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF031820.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Mar 31 14:11:44 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.83	152	278622	20.00	ng	0.00
21) Naphthalene-d8	8.12	136	977330	20.00	ng	0.00
39) Acenaphthene-d10	9.88	164	427488	20.00	ng	0.00
64) Phenanthrene-d10	11.37	188	642195	20.00	ng	0.00
76) Chrysene-d12	14.02	240	508427	20.00	ng	0.00
87) Perylene-d12	15.49	264	405917	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.45	112	1566077	89.48	ng	0.01
7) Phenol-d6	6.46	99	1955313	87.45	ng	0.00
23) Nitrobenzene-d5	7.39	82	1135181	63.13	ng	0.00
42) 2,4,6-Tribromophenol	10.67	330	339158	76.90	ng	0.00
45) 2-Fluorobiphenyl	9.20	172	1906753	66.08	ng	0.00
79) Terphenyl-d14	12.96	244	1578925	60.84	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.59	163	199067	6.556	ng	99
71) Phenanthrene	11.39	178	174499	5.240	ng	97
72) Anthracene	11.45	178	136663	4.038	ng	97
75) Fluoranthene	12.59	202	657899	18.256	ng	97
78) Pyrene	12.82	202	559897	13.418	ng	97
81) Benzo(a)anthracene	14.00	228	328762m	9.944	ng	
83) Chrysene	14.04	228	313039	9.174	ng	96
86) Indeno(1,2,3-cd)pyrene	16.95	276	113375	3.528	ng	98
88) Benzo(b)fluoranthene	15.06	252	316538m	11.674	ng	
89) Benzo(k)fluoranthene	15.08	252	104125m	4.033	ng	
90) Benzo(a)pyrene	15.42	252	204039	8.280	ng	97
92) Benzo(a,h,i)perylene	17.39	276	103994	4.803	ng	# 90

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

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