

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF050120\
 Data File : BF120273.D
 Acq On : 1 May 2020 22:08
 Operator : MA/SJ
 Sample : L2463-20
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 JHDS-SOUTH-C

Manual Integrations
 APPROVED

mohammad
 5/4/2020 3:45:39 PM

Quant Time: May 02 02:38:21 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF040820.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri May 01 15:26:36 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.65	152	90317	20.00	ng	0.00
21) Naphthalene-d8	7.93	136	324596	20.00	ng	0.00
39) Acenaphthene-d10	9.69	164	146498	20.00	ng	-0.01
64) Phenanthrene-d10	11.17	188	255565	20.00	ng	0.00
76) Chrysene-d12	13.82	240	220766	20.00	ng	-0.01
87) Perylene-d12	15.23	264	217926	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.26	112	540096	103.35	ng	0.00
7) Phenol-d6	6.31	99	672174	99.82	ng	0.00
23) Nitrobenzene-d5	7.22	82	409726	65.30	ng	-0.01
42) 2,4,6-Tribromophenol	10.48	330	136304	75.99	ng	-0.01
45) 2-Fluorobiphenyl	9.02	172	775479	75.16	ng	0.00
79) Terphenyl-d14	12.77	244	757120	57.71	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.41	163	101118	8.924	ng	98
71) Phenanthrene	11.20	178	49933	3.671	ng	98
75) Fluoranthene	12.39	202	107478	7.324	ng	94
78) Pyrene	12.62	202	99779	5.361	ng	98
81) Benzo(a)anthracene	13.80	228	47961m	3.302	ng	
83) Chrysene	13.84	228	51909	3.518	ng	95
86) Indeno(1,2,3-cd)pyrene	16.57	276	27781	2.136	ng	95
88) Benzo(b)fluoranthene	14.83	252	65986m	4.209	ng	
90) Benzo(a)pyrene	15.17	252	32958	2.500	ng	# 76
92) Benzo(g,h,i)perylene	16.98	276	31354	2.720	ng	# 85

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

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