

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG060120\
 Data File : BG045565.D
 Acq On : 1 Jun 2020 21:17
 Operator : CG/JU
 Sample : L2798-10
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 NM100-COMP-2020-0-6

Manual Integrations
 APPROVED

mohammad
 6/2/2020 3:23:07 PM

Quant Time: Jun 02 01:52:16 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG051520.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 01 15:07:00 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.95	152	79803	20.00	ng	0.00
21) Naphthalene-d8	10.76	136	328648	20.00	ng	0.00
39) Acenaphthene-d10	14.60	164	218095	20.00	ng	0.00
64) Phenanthrene-d10	17.35	188	454876	20.00	ng	0.00
76) Chrysene-d12	21.61	240	397551	20.00	ng	0.00
87) Perylene-d12	24.76	264	428057	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.55	112	353619	86.60	ng	0.00
7) Phenol-d6	7.15	99	522719	89.94	ng	0.00
23) Nitrobenzene-d5	9.11	82	361969	60.06	ng	0.00
42) 2,4,6-Tribromophenol	16.09	330	229854	82.41	ng	0.00
45) 2-Fluorobiphenyl	13.22	172	833168	64.80	ng	0.00
79) Terphenyl-d14	19.96	244	1107340	64.96	ng	0.00
Target Compounds						
50) Dimethylphthalate	14.05	163	166242	11.112	ng	98
71) Phenanthrene	17.39	178	270771	13.635	ng	97
72) Anthracene	17.48	178	62053	3.112	ng	98
75) Fluoranthene	19.40	202	534906	21.178	ng	98
78) Pyrene	19.76	202	466926	20.604	ng	98
81) Benzo(a)anthracene	21.58	228	265853	12.004	ng	99
83) Chrysene	21.65	228	247417	11.619	ng	98
86) Indeno(1,2,3-cd)pyrene	28.32	276	179538	6.447	ng	99
88) Benzo(b)fluoranthene	23.76	252	360549	15.705	ng	100
89) Benzo(k)fluoranthene	23.82	252	117603m	5.453	ng	
90) Benzo(a)pyrene	24.60	252	255514	11.951	ng	98
91) Dibenzo(a,h)anthracene	28.36	278	43473	2.023	ng	94
92) Benzo(a,h,i)perylene	29.44	276	160276	7.459	ng	97

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

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