

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG060120\
 Data File : BG045564.D
 Acq On : 1 Jun 2020 20:36
 Operator : CG/JU
 Sample : L2798-07
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 NM57-COMP-2020-0-6

Manual Integrations
 APPROVED

mohammad
 6/2/2020 3:22:59 PM

Quant Time: Jun 02 01:44:29 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	85551	20.00	ng	0.00
21) Naphthalene-d8	10.76	136	342236	20.00	ng	0.00
39) Acenaphthene-d10	14.60	164	224954	20.00	ng	0.00
64) Phenanthrene-d10	17.34	188	471201	20.00	ng	0.00
76) Chrysene-d12	21.60	240	428310	20.00	ng	0.00
87) Perylene-d12	24.75	264	459798	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.54	112	322573	73.69	ng	0.00
7) Phenol-d6	7.15	99	494751	79.41	ng	0.00
23) Nitrobenzene-d5	9.11	82	356523	56.81	ng	0.00
42) 2,4,6-Tribromophenol	16.09	330	225072	78.23	ng	0.00
45) 2-Fluorobiphenyl	13.21	172	815775	61.51	ng	0.00
79) Terphenyl-d14	19.96	244	1112256	60.57	ng	0.00
Target Compounds						
50) Dimethylphthalate	14.05	163	157240	10.189	ng	99
71) Phenanthrene	17.39	178	67515	3.282	ng	93
75) Fluoranthene	19.40	202	175529	6.709	ng	99
78) Pyrene	19.76	202	164963	6.757	ng	97
81) Benzo(a)anthracene	21.59	228	93917	3.936	ng	97
83) Chrysene	21.65	228	93412m	4.072	ng	
86) Indeno(1,2,3-cd)pyrene	28.32	276	73035	2.434	ng	# 94
88) Benzo(b)fluoranthene	23.76	252	136291	5.527	ng	# 95
89) Benzo(k)fluoranthene	23.81	252	54671m	2.360	ng	
90) Benzo(a)pyrene	24.60	252	101248	4.409	ng	98
92) Benzo(g,h,i)perylene	29.43	276	71204	3.085	ng	# 94

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

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