

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG060120\
 Data File : BG045562.D
 Acq On : 1 Jun 2020 19:16
 Operator : CG/JU
 Sample : L2773-12
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 NM48-COMP-2020-0-6

Manual Integrations
 APPROVED

mohammad
 6/2/2020 3:22:49 PM

Quant Time: Jun 02 01:26:59 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	82016	20.00	ng	0.00
21) Naphthalene-d8	10.76	136	341985	20.00	ng	0.00
39) Acenaphthene-d10	14.60	164	232879	20.00	ng	0.00
64) Phenanthrene-d10	17.35	188	493043	20.00	ng	0.00
76) Chrysene-d12	21.60	240	441424	20.00	ng	0.00
87) Perylene-d12	24.75	264	468185	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.54	112	344138	82.00	ng	0.00
7) Phenol-d6	7.15	99	523094	87.57	ng	0.00
23) Nitrobenzene-d5	9.11	82	370355	59.06	ng	0.00
42) 2,4,6-Tribromophenol	16.09	330	244056	81.95	ng	0.00
45) 2-Fluorobiphenyl	13.22	172	844357	61.50	ng	0.00
79) Terphenyl-d14	19.96	244	1176435	62.16	ng	0.00
Target Compounds						
50) Dimethylphthalate	14.04	163	158141	9.899	ng	97
71) Phenanthrene	17.39	178	123250	5.726	ng	96
75) Fluoranthene	19.40	202	314714	11.495	ng	98
78) Pyrene	19.76	202	299999	11.922	ng	98
81) Benzo(a)anthracene	21.58	228	181492	7.381	ng	99
83) Chrysene	21.65	228	171417	7.250	ng	96
86) Indeno(1,2,3-cd)pyrene	28.31	276	145620	4.710	ng	# 77
88) Benzo(b)fluoranthene	23.75	252	282312	11.243	ng	100
89) Benzo(k)fluoranthene	23.81	252	84059m	3.563	ng	
90) Benzo(a)pyrene	24.59	252	195470	8.359	ng	97
92) Benzo(g,h,i)perylene	29.43	276	149449	6.359	ng	98

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

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