

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG060120\
 Data File : BG045566.D
 Acq On : 1 Jun 2020 21:57
 Operator : CG/JU
 Sample : L2792-08
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 NM114-COMP-2020-0-6

Manual Integrations
 APPROVED

mohammad
 6/2/2020 3:23:01 PM

Quant Time: Jun 02 02:00:35 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	82324	20.00	ng	0.00
21) Naphthalene-d8	10.76	136	336163	20.00	ng	0.00
39) Acenaphthene-d10	14.60	164	219953	20.00	ng	0.00
64) Phenanthrene-d10	17.35	188	453575	20.00	ng	0.00
76) Chrysene-d12	21.60	240	393060	20.00	ng	0.00
87) Perylene-d12	24.75	264	424348	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.54	112	386418	91.73	ng	0.00
7) Phenol-d6	7.15	99	582788	97.20	ng	0.00
23) Nitrobenzene-d5	9.12	82	406230	65.90	ng	0.00
42) 2,4,6-Tribromophenol	16.09	330	261884	93.10	ng	0.00
45) 2-Fluorobiphenyl	13.22	172	917640	70.77	ng	0.00
79) Terphenyl-d14	19.96	244	1224168	72.64	ng	0.00
Target Compounds						
50) Dimethylphthalate	14.04	163	160779	10.656	ng	99
71) Phenanthrene	17.39	178	74648	3.770	ng	98
75) Fluoranthene	19.40	202	208931	8.296	ng	97
78) Pyrene	19.76	202	195277	8.715	ng	98
81) Benzo(a)anthracene	21.59	228	116738	5.331	ng	97
83) Chrysene	21.65	228	108873m	5.171	ng	
86) Indeno(1,2,3-cd)pyrene	28.32	276	76225	2.769	ng	# 93
88) Benzo(b)fluoranthene	23.76	252	183862m	8.079	ng	
90) Benzo(a)pyrene	24.61	252	116526	5.498	ng	97
92) Benzo(g,h,i)perylene	29.44	276	72597	3.408	ng	95

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

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