

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG060120\
 Data File : BG045560.D
 Acq On : 1 Jun 2020 17:56
 Operator : CG/JU
 Sample : L2792-12
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 NM75-COMP-2020-0-6

Manual Integrations
 APPROVED

mohammad
 6/2/2020 3:22:44 PM

Quant Time: Jun 01 18:43:40 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	66341	20.00	ng	0.00
21) Naphthalene-d8	10.76	136	277589	20.00	ng	0.00
39) Acenaphthene-d10	14.60	164	204522	20.00	ng	0.00
64) Phenanthrene-d10	17.34	188	439241	20.00	ng	0.00
76) Chrysene-d12	21.60	240	399928	20.00	ng	0.00
87) Perylene-d12	24.75	264	436379	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.54	112	311230	91.68	ng	0.00
7) Phenol-d6	7.15	99	492583	101.95	ng	0.00
23) Nitrobenzene-d5	9.11	82	371956	73.07	ng	0.00
42) 2,4,6-Tribromophenol	16.09	330	259054	99.04	ng	0.00
45) 2-Fluorobiphenyl	13.22	172	898439	74.51	ng	0.00
79) Terphenyl-d14	19.96	244	1326322	77.35	ng	0.00
Target Compounds						
50) Dimethylphthalate	14.05	163	107995	7.697	ng	98
71) Phenanthrene	17.39	178	91780	4.786	ng	96
75) Fluoranthene	19.40	202	236193	9.684	ng	100
78) Pyrene	19.76	202	227962	9.999	ng	97
81) Benzo(a)anthracene	21.59	228	139616	6.267	ng	99
83) Chrysene	21.65	228	132845	6.201	ng	99
86) Indeno(1,2,3-cd)pyrene	28.31	276	102659	3.665	ng	# 95
88) Benzo(b)fluoranthene	23.76	252	195778	8.365	ng	99
89) Benzo(k)fluoranthene	23.81	252	75945m	3.454	ng	
90) Benzo(a)pyrene	24.60	252	142974	6.560	ng	99
92) Benzo(g,h,i)perylene	29.42	276	103125	4.708	ng	95

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

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