

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG060120\
 Data File : BG045563.D
 Acq On : 1 Jun 2020 19:56
 Operator : CG/JU
 Sample : L2798-06
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 NM106-COMP-2020-0-6

Manual Integrations
 APPROVED

mohammad
 6/2/2020 3:22:56 PM

Quant Time: Jun 02 01:36:11 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	7.96	152	80534	20.00	ng	0.00
21) Naphthalene-d8	10.76	136	326221	20.00	ng	0.00
39) Acenaphthene-d10	14.59	164	221153	20.00	ng	0.00
64) Phenanthrene-d10	17.34	188	463132	20.00	ng	0.00
76) Chrysene-d12	21.60	240	401104	20.00	ng	0.00
87) Perylene-d12	24.76	264	430614	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.54	112	318012	77.17	ng	0.00
7) Phenol-d6	7.15	99	478312	81.55	ng	0.00
23) Nitrobenzene-d5	9.11	82	340994	57.00	ng	0.00
42) 2,4,6-Tribromophenol	16.09	330	220373	77.92	ng	0.00
45) 2-Fluorobiphenyl	13.22	172	781091	59.91	ng	0.00
79) Terphenyl-d14	19.96	244	1084011	63.03	ng	0.00
Target Compounds						
50) Dimethylphthalate	14.05	163	150191	9.900	ng	99
71) Phenanthrene	17.38	178	122563	6.062	ng	94
75) Fluoranthene	19.40	202	298252	11.598	ng	100
78) Pyrene	19.76	202	278167	12.166	ng	98
81) Benzo(a)anthracene	21.58	228	157579	7.052	ng	99
83) Chrysene	21.65	228	141586	6.590	ng	98
86) Indeno(1,2,3-cd)pyrene	28.32	276	110097	3.919	ng	# 97
88) Benzo(b)fluoranthene	23.76	252	223061	9.659	ng	98
89) Benzo(k)fluoranthene	23.82	252	65601m	3.024	ng	
90) Benzo(a)pyrene	24.60	252	156110	7.258	ng	# 94
92) Benzo(g,h,i)perylene	29.43	276	103159	4.772	ng	93

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

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