

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG060120\
 Data File : BG045568.D
 Acq On : 1 Jun 2020 23:18
 Operator : CG/JU
 Sample : L2798-03
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 NM87-COMP-2020-0-6

Manual Integrations
 APPROVED

mohammad
 6/2/2020 3:23:10 PM

Quant Time: Jun 02 02:13:47 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	81279	20.00	ng	0.00
21) Naphthalene-d8	10.76	136	318724	20.00	ng	0.00
39) Acenaphthene-d10	14.60	164	212257	20.00	ng	0.00
64) Phenanthrene-d10	17.35	188	442163	20.00	ng	0.00
76) Chrysene-d12	21.60	240	376222	20.00	ng	0.00
87) Perylene-d12	24.76	264	417869	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.54	112	294789	70.88	ng	0.00
7) Phenol-d6	7.15	99	490703	82.90	ng	0.00
23) Nitrobenzene-d5	9.12	82	327611	56.05	ng	0.00
42) 2,4,6-Tribromophenol	16.09	330	232690	85.72	ng	0.00
45) 2-Fluorobiphenyl	13.22	172	791202	63.23	ng	0.00
79) Terphenyl-d14	19.96	244	1064611	66.00	ng	0.00
Target Compounds						
50) Dimethylphthalate	14.05	163	155035	10.648	ng	99
71) Phenanthrene	17.39	178	102494	5.310	ng	94
75) Fluoranthene	19.40	202	235069	9.574	ng	99
78) Pyrene	19.76	202	227902	10.627	ng	95
81) Benzo(a)anthracene	21.59	228	130966	6.249	ng	96
83) Chrysene	21.65	228	123992	6.153	ng	99
86) Indeno(1,2,3-cd)pyrene	28.33	276	83754	3.178	ng	# 91
88) Benzo(b)fluoranthene	23.76	252	183536	8.190	ng	# 95
89) Benzo(k)fluoranthene	23.82	252	68600m	3.258	ng	
90) Benzo(a)pyrene	24.61	252	131258	6.289	ng	95
92) Benzo(g,h,i)perylene	29.44	276	77778	3.708	ng	97

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

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