

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF060220\
 Data File : BF120497.D
 Acq On : 2 Jun 2020 16:17
 Operator : JU/CG
 Sample : L2798-11
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 NM101-COMP-2020-0-6

Manual Integrations
 APPROVED

mohammad
 6/3/2020 4:36:12 PM

Quant Time: Jun 03 03:56:40 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF052820.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 02 11:37:51 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.85	152	212070	20.00	ng	0.00
21) Naphthalene-d8	8.13	136	760440	20.00	ng	0.00
39) Acenaphthene-d10	9.88	164	361901	20.00	ng	0.00
64) Phenanthrene-d10	11.37	188	586426	20.00	ng	0.00
76) Chrysene-d12	14.01	240	411496	20.00	ng	0.00
86) Perylene-d12	15.47	264	395664	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.47	112	941278	84.59	ng	0.00
7) Phenol-d6	6.48	99	1189441	86.72	ng	0.00
23) Nitrobenzene-d5	7.40	82	774415	66.62	ng	0.00
42) 2,4,6-Tribromophenol	10.67	330	294584	87.85	ng	0.00
45) 2-Fluorobiphenyl	9.20	172	1414192	66.30	ng	0.00
79) Terphenyl-d14	12.96	244	1198947	64.84	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.59	163	268266	11.857	ng	99
71) Phenanthrene	11.39	178	310444	12.014	ng	98
72) Anthracene	11.44	178	82822	3.195	ng	98
75) Fluoranthene	12.58	202	571173	20.543	ng	98
78) Pyrene	12.81	202	496973	17.366	ng	98
81) Benzo(a)anthracene	14.00	228	284218	12.887	ng	98
83) Chrysene	14.03	228	247769	10.716	ng	97
87) Indeno(1,2,3-cd)pyrene	16.93	276	123164	5.691	ng	97
88) Benzo(b)fluoranthene	15.04	252	323026m	14.699	ng	
89) Benzo(k)fluoranthene	15.07	252	111834m	5.451	ng	
90) Benzo(a)pyrene	15.41	252	219582	11.423	ng	99
92) Benzo(a,h,i)perylene	17.37	276	115873	6.659	ng	98

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

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