

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF060220\
 Data File : BF120502.D
 Acq On : 2 Jun 2020 18:35
 Operator : JU/CG
 Sample : L2798-08
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 NM98-COMP-2020-0-6

Manual Integrations
 APPROVED

mohammad
 6/3/2020 4:36:28 PM

Quant Time: Jun 02 19:34:49 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	176454	20.00	ng	0.00
21) Naphthalene-d8	8.13	136	601451	20.00	ng	0.00
39) Acenaphthene-d10	9.88	164	268019	20.00	ng	0.00
64) Phenanthrene-d10	11.37	188	461163	20.00	ng	0.00
76) Chrysene-d12	14.01	240	368118	20.00	ng	0.00
86) Perylene-d12	15.47	264	317484	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.47	112	667919	72.14	ng	0.00
7) Phenol-d6	6.48	99	825912	72.37	ng	0.00
23) Nitrobenzene-d5	7.40	82	576646	62.72	ng	0.00
42) 2,4,6-Tribromophenol	10.67	330	203395	81.90	ng	-0.01
45) 2-Fluorobiphenyl	9.20	172	1066208	67.50	ng	0.00
79) Terphenyl-d14	12.96	244	1053612	63.69	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.59	163	190693	11.381	ng	99
71) Phenanthrene	11.39	178	160577	7.902	ng	98
72) Anthracene	11.44	178	50020	2.453	ng	97
75) Fluoranthene	12.58	202	377040	17.244	ng	99
78) Pyrene	12.81	202	371744	14.521	ng	98
81) Benzo(a)anthracene	14.00	228	199293	10.101	ng	97
83) Chrysene	14.04	228	188318	9.105	ng	98
87) Indeno(1,2,3-cd)pyrene	16.93	276	103131	5.939	ng	95
88) Benzo(b)fluoranthene	15.04	252	255584m	14.494	ng	
89) Benzo(k)fluoranthene	15.07	252	59614m	3.621	ng	
90) Benzo(a)pyrene	15.41	252	158796	10.295	ng	98
92) Benzo(a,h,i)perylene	17.37	276	97295	6.968	ng	97

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

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