

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF061820\
 Data File : BF120657.D
 Acq On : 18 Jun 2020 22:22
 Operator : JU/CG
 Sample : L2818-01 2X
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TR-05-061720

Manual Integrations
 APPROVED

mohammad
 6/19/2020 12:58:19 PM

Quant Time: Jun 19 02:06:49 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jun 11 11:28:22 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.80	152	139085	20.00	ng	0.00
21) Naphthalene-d8	8.09	136	525966	20.00	ng	0.00
39) Acenaphthene-d10	9.84	164	266767	20.00	ng	0.00
64) Phenanthrene-d10	11.32	188	507260	20.00	ng	0.00
76) Chrysene-d12	13.96	240	417430	20.00	ng	0.00
86) Perylene-d12	15.41	264	357935	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.42	112	321678	39.65	ng	0.00
7) Phenol-d6	6.44	99	392145	38.81	ng	0.00
23) Nitrobenzene-d5	7.36	82	216512	23.01	ng	0.00
42) 2,4,6-Tribromophenol	10.63	330	118553	38.22	ng	0.00
45) 2-Fluorobiphenyl	9.16	172	504243	28.92	ng	0.00
79) Terphenyl-d14	12.90	244	610074	32.28	ng	0.00
Target Compounds						Qvalue
71) Phenanthrene	11.35	178	96554	3.829	ng	98
75) Fluoranthene	12.53	202	268084	9.299	ng	99
78) Pyrene	12.76	202	219542	7.424	ng	98
81) Benzo(a)anthracene	13.95	228	107738	4.296	ng	100
83) Chrysene	13.99	228	126977	4.936	ng	96
87) Indeno(1,2,3-cd)pyrene	16.83	276	51127	2.348	ng	97
88) Benzo(b)fluoranthene	14.99	252	154225m	7.233	ng	
90) Benzo(a)pyrene	15.34	252	78286	4.103	ng	# 94
92) Benzo(g,h,i)perylene	17.26	276	48626	2.763	ng	# 89

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

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