

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF061820\
 Data File : BF120656.D
 Acq On : 18 Jun 2020 21:54
 Operator : JU/CG
 Sample : L2813-07 2X
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 EO-01-061720

Manual Integrations
 APPROVED

mohammad
 6/19/2020 12:58:16 PM

Quant Time: Jun 19 02:02:57 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	139597	20.00	ng	0.00
21) Naphthalene-d8	8.09	136	524573	20.00	ng	0.00
39) Acenaphthene-d10	9.84	164	266786	20.00	ng	0.00
64) Phenanthrene-d10	11.32	188	506739	20.00	ng	0.00
76) Chrysene-d12	13.96	240	409492	20.00	ng	0.00
86) Perylene-d12	15.41	264	358020	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.42	112	225464	27.69	ng	0.00
7) Phenol-d6	6.44	99	277896	27.40	ng	0.00
23) Nitrobenzene-d5	7.36	82	159036	16.95	ng	0.00
42) 2,4,6-Tribromophenol	10.63	330	82725	26.67	ng	0.00
45) 2-Fluorobiphenyl	9.16	172	378783	21.72	ng	0.00
79) Terphenyl-d14	12.90	244	458804	24.75	ng	0.00
Target Compounds						
71) Phenanthrene	11.35	178	64111	2.545	ng	97
75) Fluoranthene	12.53	202	111547	3.873	ng	100
78) Pyrene	12.76	202	113709	3.920	ng	97
81) Benzo(a)anthracene	13.95	228	55632	2.261	ng	# 94
83) Chrysene	13.99	228	55199	2.187	ng	# 93
88) Benzo(b)fluoranthene	14.99	252	71713m	3.363	ng	
90) Benzo(a)pyrene	15.34	252	48505	2.541	ng	# 83

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

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