

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062920\
 Data File : BF120737.D
 Acq On : 29 Jun 2020 19:46
 Operator : JU/CG
 Sample : L2819-06
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 BU-03-062620

Manual Integrations
 APPROVED

mohammad
 6/30/2020 4:14:41 PM

Quant Time: Jun 30 02:00:13 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 29 15:30:14 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.80	152	57182	20.00	ng	0.00
21) Naphthalene-d8	8.08	136	200306	20.00	ng	0.00
39) Acenaphthene-d10	9.83	164	106772	20.00	ng	0.00
64) Phenanthrene-d10	11.32	188	441268	20.00	ng	0.00
76) Chrysene-d12	13.96	240	146873	20.00	ng	0.00
86) Perylene-d12	15.41	264	156575	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.48	112	324545	97.31	ng	0.08
7) Phenol-d6	6.51	99	385934	92.90	ng	0.08
23) Nitrobenzene-d5	7.36	82	245295	68.45	ng	0.00
42) 2,4,6-Tribromophenol	10.63	330	123857	99.77	ng	0.01
45) 2-Fluorobiphenyl	9.15	172	533180	76.40	ng	0.00
79) Terphenyl-d14	12.91	244	554643	83.41	ng	0.01
Target Compounds						
32) Benzoic acid	7.86	122	6386	2.849	ng	95
50) Dimethylphthalate	9.55	163	52637	6.845	ng	100
78) Pyrene	12.76	202	37198	3.575	ng	96
81) Benzo(a)anthracene	13.95	228	23806	2.698	ng	96
83) Chrysene	13.99	228	22236	2.457	ng	95
88) Benzo(b)fluoranthene	14.99	252	29192m	3.130	ng	
90) Benzo(a)pyrene	15.34	252	19634	2.352	ng	# 86

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

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