

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF081919\
 Data File : BF116409.D
 Acq On : 20 Aug 2019 1:05
 Operator : HP/JU
 Sample : K4223-11
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SP-6

Manual Integrations
 APPROVED

mohammad
 8/22/2019 12:45:48 PM

Quant Time: Aug 20 03:55:34 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF073019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 19 06:03:31 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.80	152	105612	20.00	ng	-0.01
21) Naphthalene-d8	8.08	136	409714	20.00	ng	-0.02
39) Acenaphthene-d10	9.83	164	226538	20.00	ng	-0.02
64) Phenanthrene-d10	11.32	188	413353	20.00	ng	-0.02
76) Chrysene-d12	13.96	240	259915	20.00	ng	-0.02
87) Perylene-d12	15.42	264	302740	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.43	112	632488	100.26	ng	0.00
7) Phenol-d6	6.45	99	850078	106.80	ng	-0.01
23) Nitrobenzene-d5	7.37	82	528191	74.41	ng	-0.02
42) 2,4,6-Tribromophenol	10.63	330	236599	107.72	ng	-0.02
45) 2-Fluorobiphenyl	9.16	172	847520	70.08	ng	-0.01
79) Terphenyl-d14	12.90	244	855107	69.32	ng	-0.02
Target Compounds						
50) Dimethylphthalate	9.55	163	201820	13.084	ng	98
71) Phenanthrene	11.35	178	49846	2.359	ng	98
75) Fluoranthene	12.54	202	85632	3.871	ng	98
78) Pyrene	12.77	202	72358	3.693	ng	98
88) Benzo(b)fluoranthene	15.01	252	42123m	2.406	ng	

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

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