

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF122619\
 Data File : BF118315.D
 Acq On : 26 Dec 2019 14:36
 Operator : JU
 Sample : K6421-01
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-2

Manual Integrations
 APPROVED

mohammad
 12/27/2019 3:19:08 PM

Quant Time: Dec 27 04:40:37 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF122419.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Dec 24 15:58:52 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.85	152	113653	20.00	ng	0.00
21) Naphthalene-d8	8.13	136	450681	20.00	ng	0.00
39) Acenaphthene-d10	9.90	164	244648	20.00	ng	0.00
64) Phenanthrene-d10	11.39	188	471215	20.00	ng	0.00
76) Chrysene-d12	14.04	240	320467	20.00	ng	-0.01
87) Perylene-d12	15.53	264	327470	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.46	112	567359	85.21	ng	0.00
7) Phenol-d6	6.48	99	721290	87.33	ng	0.00
23) Nitrobenzene-d5	7.41	82	426318	54.10	ng	-0.01
42) 2,4,6-Tribromophenol	10.69	330	246985	81.75	ng	0.00
45) 2-Fluorobiphenyl	9.22	172	903842	56.39	ng	0.00
79) Terphenyl-d14	12.98	244	1173055	60.73	ng	0.00
Target Compounds						
10) Phenol	6.49	94	26619	2.867	ng	Qvalue # 65
50) Dimethylphthalate	9.61	163	48668	2.650	ng	97
71) Phenanthrene	11.41	178	139469	5.429	ng	98
75) Fluoranthene	12.61	202	196871	7.594	ng	99
78) Pyrene	12.84	202	169701	6.485	ng	99
81) Benzo(a)anthracene	14.03	228	87221	3.888	ng	98
83) Chrysene	14.06	228	72607	3.380	ng	97
88) Benzo(b)fluoranthene	15.09	252	81885m	3.767	ng	
90) Benzo(a)pyrene	15.46	252	61577	3.216	ng	97
92) Benzo(g,h,i)perylene	17.47	276	34123	2.126	ng	96

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

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