

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF081919\
 Data File : BF116403.D
 Acq On : 19 Aug 2019 22:22
 Operator : HP/JU
 Sample : K4223-09
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SP-5

Manual Integrations
 APPROVED

mohammad
 8/22/2019 12:45:34 PM

Quant Time: Aug 20 11:32:25 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	98101	20.00	ng	-0.01
21) Naphthalene-d8	8.08	136	388821	20.00	ng	-0.02
39) Acenaphthene-d10	9.83	164	212955	20.00	ng	-0.02
64) Phenanthrene-d10	11.32	188	403774	20.00	ng	-0.02
76) Chrysene-d12	13.96	240	295209	20.00	ng	-0.02
87) Perylene-d12	15.42	264	293951	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.43	112	610100	104.11	ng	-0.01
7) Phenol-d6	6.45	99	815116	110.25	ng	-0.02
23) Nitrobenzene-d5	7.37	82	498518	74.01	ng	-0.02
42) 2,4,6-Tribromophenol	10.63	330	224451	108.71	ng	-0.02
45) 2-Fluorobiphenyl	9.15	172	758586	66.72	ng	-0.02
79) Terphenyl-d14	12.90	244	900275	64.26	ng	-0.02
Target Compounds						
50) Dimethylphthalate	9.55	163	208867	14.405	ng	99
71) Phenanthrene	11.35	178	51191	2.480	ng	98
75) Fluoranthene	12.54	202	96193	4.451	ng	99
78) Pyrene	12.77	202	86375	3.881	ng	97
81) Benzo(a)anthracene	13.96	228	38909m	2.062	ng	
83) Chrysene	13.99	228	39453m	2.154	ng	
88) Benzo(b)fluoranthene	15.01	252	41882m	2.464	ng	

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

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