

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF090619\
 Data File : BF116856.D
 Acq On : 6 Sep 2019 12:40
 Operator : HP/JU
 Sample : K4650-05
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-3

Manual Integrations
 APPROVED

mohammad
 9/12/2019 9:11:28 AM

Quant Time: Sep 07 04:08:14 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF083019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 30 20:02:30 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.85	152	79697	20.00	ng	0.00
21) Naphthalene-d8	8.13	136	321619	20.00	ng	0.00
39) Acenaphthene-d10	9.88	164	166797	20.00	ng	0.00
64) Phenanthrene-d10	11.36	188	279535	20.00	ng	0.00
76) Chrysene-d12	14.00	240	202142	20.00	ng	0.00
87) Perylene-d12	15.49	264	189704	20.00	ng	0.05
System Monitoring Compounds						
5) 2-Fluorophenol	5.47	112	637713	129.63	ng	0.00
7) Phenol-d6	6.49	99	861862	133.42	ng	0.00
23) Nitrobenzene-d5	7.40	82	540507	95.94	ng	0.00
42) 2,4,6-Tribromophenol	10.67	330	181963	163.18	ng	0.00
45) 2-Fluorobiphenyl	9.20	172	867015	87.69	ng	0.00
79) Terphenyl-d14	12.95	244	840963	76.96	ng	0.00
Target Compounds						
10) Phenol	6.49	94	24733	3.458	ng	87
50) Dimethylphthalate	9.59	163	157894	13.711	ng	97
75) Fluoranthene	12.57	202	67469	4.412	ng	97
78) Pyrene	12.80	202	59025	3.285	ng	98
81) Benzo(a)anthracene	14.00	228	28221m	2.206	ng	
83) Chrysene	14.03	228	30107	2.284	ng	99
88) Benzo(b)fluoranthene	15.06	252	37039m	3.229	ng	
90) Benzo(a)pyrene	15.43	252	23758	2.381	ng	96

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

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