

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF090419\
 Data File : BF116810.D
 Acq On : 4 Sep 2019 12:50
 Operator : HP/JU
 Sample : K4601-01
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 T12-E1-E4-(0-2.7)

Manual Integrations
 APPROVED

mohammad
 9/9/2019 9:55:39 AM

Quant Time: Sep 05 03:18:48 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	75228	20.00	ng	0.00
21) Naphthalene-d8	8.13	136	300459	20.00	ng	0.00
39) Acenaphthene-d10	9.88	164	157660	20.00	ng	0.00
64) Phenanthrene-d10	11.36	188	251610	20.00	ng	0.00
76) Chrysene-d12	14.00	240	180544	20.00	ng	0.00
87) Perylene-d12	15.45	264	176844	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.46	112	450115	96.93	ng	0.00
7) Phenol-d6	6.47	99	620104	101.70	ng	0.00
23) Nitrobenzene-d5	7.40	82	375374	71.32	ng	0.00
42) 2,4,6-Tribromophenol	10.67	330	129311	122.69	ng	0.00
45) 2-Fluorobiphenyl	9.20	172	649081	69.45	ng	0.00
79) Terphenyl-d14	12.95	244	619017	63.43	ng	0.00
Target Compounds						
10) Phenol	6.49	94	34419	5.098	ng	96
50) Dimethylphthalate	9.59	163	84310	7.746	ng	98
75) Fluoranthene	12.58	202	52543	3.817	ng	81
78) Pyrene	12.80	202	45902	2.860	ng	# 85
83) Chrysene	14.03	228	26906	2.286	ng	# 85
88) Benzo(b)fluoranthene	15.03	252	33050m	3.091	ng	

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

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