

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF090519\
 Data File : BF116840.D
 Acq On : 5 Sep 2019 16:14
 Operator : HP/JU
 Sample : K4683-01
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 20190663-LOCATION-1

Manual Integrations
 APPROVED

mohammad
 9/10/2019 9:52:22 AM

Quant Time: Sep 06 04:40:21 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	86783	20.00	ng	0.00
21) Naphthalene-d8	8.13	136	357003	20.00	ng	0.00
39) Acenaphthene-d10	9.88	164	186230	20.00	ng	0.00
64) Phenanthrene-d10	11.36	188	297741	20.00	ng	0.00
76) Chrysene-d12	14.00	240	224005	20.00	ng	0.00
87) Perylene-d12	15.46	264	215931	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.47	112	580528	108.37	ng	0.00
7) Phenol-d6	6.48	99	794517	112.95	ng	0.00
23) Nitrobenzene-d5	7.40	82	482501	77.16	ng	0.00
42) 2,4,6-Tribromophenol	10.67	330	154925	124.44	ng	0.00
45) 2-Fluorobiphenyl	9.20	172	775252	70.22	ng	-0.01
79) Terphenyl-d14	12.95	244	738118	60.96	ng	0.00
Target Compounds						
10) Phenol	6.49	94	19670	2.525	ng	97
50) Dimethylphthalate	9.59	163	87063	6.772	ng	98
71) Phenanthrene	11.39	178	57044	3.442	ng	97
75) Fluoranthene	12.57	202	49490	3.038	ng	98
78) Pyrene	12.80	202	62030	3.115	ng	98
81) Benzo(a)anthracene	13.99	228	30536	2.154	ng	# 78
83) Chrysene	14.02	228	33650	2.304	ng	# 94
88) Benzo(b)fluoranthene	15.03	252	27528m	2.109	ng	

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

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