

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF090119\
 Data File : BF116752.D
 Acq On : 1 Sep 2019 19:36
 Operator : HP/JU
 Sample : K4547-07
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 T3-S-G

Manual Integrations
 APPROVED

mohammad
 9/5/2019 9:54:47 AM

Quant Time: Sep 02 06:39:56 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	75089	20.00	ng	0.00
21) Naphthalene-d8	8.13	136	300503	20.00	ng	0.00
39) Acenaphthene-d10	9.88	164	173374	20.00	ng	0.00
64) Phenanthrene-d10	11.36	188	314760	20.00	ng	0.00
76) Chrysene-d12	14.00	240	220931	20.00	ng	0.00
87) Perylene-d12	15.44	264	237069	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.46	112	439932	94.92	ng	0.00
7) Phenol-d6	6.47	99	604527	99.33	ng	-0.01
23) Nitrobenzene-d5	7.40	82	370934	70.47	ng	-0.01
42) 2,4,6-Tribromophenol	10.66	330	139004	119.93	ng	-0.01
45) 2-Fluorobiphenyl	9.20	172	674650	65.64	ng	-0.01
79) Terphenyl-d14	12.94	244	751910	62.96	ng	0.00
Target Compounds						
10) Phenol	6.48	94	30807	4.571	ng	96
50) Dimethylphthalate	9.59	163	67855	5.669	ng	98
71) Phenanthrene	11.39	178	44579	2.544	ng	# 91
75) Fluoranthene	12.57	202	129743	7.535	ng	94
78) Pyrene	12.80	202	110967	5.650	ng	# 93
81) Benzo(a)anthracene	13.99	228	51586	3.689	ng	95
83) Chrysene	14.02	228	69710	4.839	ng	95
84) Bis(2-ethylhexyl)phthalate	13.98	149	98269	8.335	ng	# 98
86) Indeno(1,2,3-cd)pyrene	16.88	276	42581	3.680	ng	96
88) Benzo(b)fluoranthene	15.02	252	104766m	7.310	ng	
89) Benzo(k)fluoranthene	15.04	252	35743m	2.735	ng	
90) Benzo(a)pyrene	15.38	252	61920	4.966	ng	# 91
92) Benzo(a,h,i)perylene	17.31	276	41073	3.690	ng	# 90

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

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