

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF090119\
 Data File : BF116754.D
 Acq On : 1 Sep 2019 20:30
 Operator : HP/JU
 Sample : K4547-15
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 T7-S-G1-G4

Manual Integrations
 APPROVED

mohammad
 9/5/2019 9:54:53 AM

Quant Time: Sep 02 06:46:53 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	91979	20.00	ng	0.00
21) Naphthalene-d8	8.13	136	384323	20.00	ng	0.00
39) Acenaphthene-d10	9.88	164	215806	20.00	ng	0.00
64) Phenanthrene-d10	11.36	188	353378	20.00	ng	0.00
76) Chrysene-d12	13.99	240	218662	20.00	ng	0.00
87) Perylene-d12	15.45	264	211896	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.46	112	497154	87.57	ng	0.00
7) Phenol-d6	6.47	99	685215	91.91	ng	0.00
23) Nitrobenzene-d5	7.40	82	410744	61.01	ng	0.00
42) 2,4,6-Tribromophenol	10.67	330	134870	93.48	ng	0.00
45) 2-Fluorobiphenyl	9.20	172	776163	60.67	ng	0.00
79) Terphenyl-d14	12.94	244	707596	59.86	ng	0.00
Target Compounds						
10) Phenol	6.49	94	52958	6.415	ng	95
50) Dimethylphthalate	9.59	163	123446	8.285	ng	99
75) Fluoranthene	12.57	202	84193	4.355	ng	94
78) Pyrene	12.80	202	75201	3.869	ng	97
81) Benzo(a)anthracene	13.98	228	35923	2.596	ng	# 94
83) Chrysene	14.02	228	43437	3.047	ng	# 90
84) Bis(2-ethylhexyl)phthalate	13.97	149	24404	2.091	ng	# 99
88) Benzo(b)fluoranthene	15.03	252	59926m	4.678	ng	
90) Benzo(a)pyrene	15.39	252	33732	3.026	ng	# 73
92) Benzo(g,h,i)perylene	17.32	276	20477	2.058	ng	# 80

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

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