

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF090419\
 Data File : BF116811.D
 Acq On : 4 Sep 2019 13:17
 Operator : HP/JU
 Sample : K4624-06
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 T-11-A1-A4-(0-4)

Manual Integrations
 APPROVED

mohammad
 9/9/2019 9:55:43 AM

Quant Time: Sep 05 03:22:11 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	66993	20.00	ng	0.00
21) Naphthalene-d8	8.13	136	265497	20.00	ng	0.00
39) Acenaphthene-d10	9.88	164	134459	20.00	ng	0.00
64) Phenanthrene-d10	11.36	188	211233	20.00	ng	0.00
76) Chrysene-d12	14.02	240	137198	20.00	ng	0.02
87) Perylene-d12	15.46	264	161253	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.46	112	362736	87.72	ng	0.00
7) Phenol-d6	6.47	99	518706	95.53	ng	0.00
23) Nitrobenzene-d5	7.40	82	321728	69.18	ng	0.00
42) 2,4,6-Tribromophenol	10.67	330	113253	125.99	ng	0.00
45) 2-Fluorobiphenyl	9.20	172	569180	71.41	ng	0.00
79) Terphenyl-d14	12.96	244	468102	63.12	ng	0.01
Target Compounds						
10) Phenol	6.49	94	24948	4.149	ng	96
50) Dimethylphthalate	9.59	163	133800	14.413	ng	98
75) Fluoranthene	12.59	202	44372	3.840	ng	# 58
78) Pyrene	12.82	202	35912	2.945	ng	# 55
83) Chrysene	14.04	228	23538	2.631	ng	# 79
84) Bis(2-ethylhexyl)phthalate	13.99	149	55903	7.636	ng	# 99
86) Indeno(1,2,3-cd)pyrene	16.90	276	17306	2.408	ng	# 93
88) Benzo(b)fluoranthene	15.04	252	32205m	3.303	ng	
90) Benzo(a)pyrene	15.40	252	17469	2.060	ng	# 64
92) Benzo(g,h,i)perylene	17.33	276	15884	2.098	ng	# 71

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

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