

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF011719\
 Data File : BF112076.D
 Acq On : 17 Jan 2019 9:05
 Operator : JU/SJ
 Sample : K1147-01
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-1

Manual Integrations
 APPROVED

Sohil
 1/17/2019 2:55:35 PM

Quant Time: Jan 17 10:22:12 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF011619.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jan 16 18:22:05 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.86	152	314172	20.00	ng	0.00
21) Naphthalene-d8	8.14	136	1208701	20.00	ng	0.00
39) Acenaphthene-d10	9.90	164	589821	20.00	ng	0.00
64) Phenanthrene-d10	11.39	188	1079502	20.00	ng	0.00
76) Chrysene-d12	14.03	240	807850	20.00	ng	0.00
87) Perylene-d12	15.50	264	759930	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.50	112	1580686	91.32	ng	0.00
7) Phenol-d6	6.51	99	2000366	92.37	ng	0.00
23) Nitrobenzene-d5	7.42	82	1232420	70.44	ng	0.00
42) 2,4,6-Tribromophenol	10.69	330	610561	95.96	ng	0.00
45) 2-Fluorobiphenyl	9.22	172	2202710	54.75	ng	0.00
79) Terphenyl-d14	12.97	244	2175994	57.30	ng	0.00
Target Compounds						
10) Phenol	6.52	94	116703	4.931	ng	87
50) Dimethylphthalate	9.60	163	283339	6.587	ng	# 96
71) Phenanthrene	11.41	178	174666	3.215	ng	# 92
75) Fluoranthene	12.60	202	541248	9.436	ng	99
78) Pyrene	12.83	202	488810	9.171	ng	98
81) Benzo(a)anthracene	14.02	228	326428	7.047	ng	98
83) Chrysene	14.05	228	300572	6.858	ng	95
86) Indeno(1,2,3-cd)pyrene	16.97	276	137228	3.394	ng	95
88) Benzo(b)fluoranthene	15.07	252	376080m	8.718	ng	
89) Benzo(k)fluoranthene	15.09	252	114592m	2.745	ng	
90) Benzo(a)pyrene	15.43	252	249271	6.301	ng	97
92) Benzo(a,h,i)perylene	17.42	276	125209	3.617	ng	100

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

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