

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF011719\
 Data File : BF112089.D
 Acq On : 17 Jan 2019 15:11
 Operator : JU/SJ
 Sample : K1167-05
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 D-9

Manual Integrations
 APPROVED

Sohil
 1/18/2019 11:45:47 AM

Quant Time: Jan 18 00:27:44 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	330716	20.00	ng	0.00
21) Naphthalene-d8	8.15	136	1211592	20.00	ng	0.00
39) Acenaphthene-d10	9.90	164	567562	20.00	ng	0.00
64) Phenanthrene-d10	11.39	188	949869	20.00	ng	0.00
76) Chrysene-d12	14.03	240	768571	20.00	ng	0.00
87) Perylene-d12	15.50	264	679489	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.50	112	2281083	125.19	ng	0.01
7) Phenol-d6	6.51	99	2935071	128.75	ng	0.00
23) Nitrobenzene-d5	7.43	82	1808899	103.15	ng	0.00
42) 2,4,6-Tribromophenol	10.69	330	801309	130.87	ng	0.00
45) 2-Fluorobiphenyl	9.22	172	3147877	81.32	ng	0.00
79) Terphenyl-d14	12.97	244	2854810	79.02	ng	0.00
Target Compounds						
10) Phenol	6.52	94	197828	7.940	ng	Qvalue # 74
50) Dimethylphthalate	9.60	163	422061	10.197	ng	100
71) Phenanthrene	11.41	178	259240	5.423	ng	97
75) Fluoranthene	12.60	202	456036	9.035	ng	99
78) Pyrene	12.83	202	418333	8.250	ng	98
81) Benzo(a)anthracene	14.02	228	200044	4.539	ng	99
83) Chrysene	14.05	228	179446	4.303	ng	97
86) Indeno(1,2,3-cd)pyrene	16.98	276	87408	2.272	ng	95
88) Benzo(b)fluoranthene	15.07	252	233103m	6.044	ng	
89) Benzo(k)fluoranthene	15.09	252	79181m	2.121	ng	
90) Benzo(a)pyrene	15.44	252	148926	4.210	ng	99
92) Benzo(a,h,i)perylene	17.42	276	85561	2.764	ng	97

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

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