

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF010819\
 Data File : BF111935.D
 Acq On : 9 Jan 2019 4:56
 Operator : JU/SJ
 Sample : K1075-17DL 10X
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 BASIN-9DL

Manual Integrations
 APPROVED

Sohil
 1/9/2019 10:37:03 AM

Quant Time: Jan 09 05:35:19 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF010719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jan 07 14:33:51 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.89	152	100602	20.00	ng	0.00
21) Naphthalene-d8	8.17	136	321084	20.00	ng	0.00
39) Acenaphthene-d10	9.93	164	145881	20.00	ng	0.00
64) Phenanthrene-d10	11.42	188	296975	20.00	ng	0.00
76) Chrysene-d12	14.06	240	241511	20.00	ng	0.00
87) Perylene-d12	15.54	264	165597	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.52	112	74304	12.64	ng	0.00
7) Phenol-d6	6.52	99	91440	12.06	ng	-0.01
23) Nitrobenzene-d5	7.44	82	46746	7.78	ng	-0.02
42) 2,4,6-Tribromophenol	10.72	330	17099	9.01	ng	0.00
45) 2-Fluorobiphenyl	9.24	172	73115	7.31	ng	-0.01
79) Terphenyl-d14	12.99	244	78668	6.18	ng	0.00
Target Compounds						
52) Acenaphthene	9.96	154	27274	2.988	ng	98
58) Fluorene	10.47	166	36913	3.821	ng	97
71) Phenanthrene	11.44	178	403329	25.402	ng	98
72) Anthracene	11.49	178	137696	8.543	ng	97
75) Fluoranthene	12.63	202	512319	31.431	ng	98
78) Pyrene	12.86	202	389519	23.440	ng	99
81) Benzo(a)anthracene	14.04	228	216356	15.566	ng	98
83) Chrysene	14.08	228	166160	12.538	ng	97
86) Indeno(1,2,3-cd)pyrene	17.03	276	43601	4.513	ng	97
88) Benzo(b)fluoranthene	15.10	252	147770m	14.917	ng	
89) Benzo(k)fluoranthene	15.13	252	56786m	6.099	ng	
90) Benzo(a)pyrene	15.47	252	102149	11.728	ng	# 97
92) Benzo(a,h,i)perylene	17.49	276	37694	5.501	ng	# 89

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

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