

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF021319\
 Data File : BF112558.D
 Acq On : 14 Feb 2019 2:45
 Operator : JU/SJ
 Sample : K1343-15DL 10X
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 JC-02-021119-HDL

Manual Integrations
 APPROVED

Sohil
 2/14/2019 11:36:29 AM

Quant Time: Feb 14 07:04:25 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF012519.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jan 28 10:38:18 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.83	152	95252	20.00	ng	-0.01
21) Naphthalene-d8	8.12	136	344470	20.00	ng	-0.01
39) Acenaphthene-d10	9.87	164	131627	20.00	ng	-0.01
64) Phenanthrene-d10	11.36	188	212048	20.00	ng	0.00
76) Chrysene-d12	14.02	240	182078	20.00	ng	0.00
87) Perylene-d12	15.49	264	125198	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.48	112	56420	12.58	ng	0.00
7) Phenol-d6	6.50	99	76950	12.35	ng	0.00
23) Nitrobenzene-d5	7.40	82	43359	7.25	ng	0.00
42) 2,4,6-Tribromophenol	10.67	330	13501	8.81	ng	0.00
45) 2-Fluorobiphenyl	9.19	172	90255	9.70	ng	-0.01
79) Terphenyl-d14	12.94	244	79091	8.18	ng	0.00
Target Compounds						
49) Acenaphthylene	9.74	152	30622	2.341	ng	98
55) Dibenzofuran	10.08	168	27896	2.336	ng	97
71) Phenanthrene	11.39	178	309786	28.101	ng	97
72) Anthracene	11.45	178	41693	3.797	ng	99
75) Fluoranthene	12.58	202	385164	32.575	ng	98
78) Pyrene	12.82	202	330809	26.242	ng	99
81) Benzo(a)anthracene	14.00	228	93883	8.771	ng	99
83) Chrysene	14.04	228	108372	10.607	ng	96
86) Indeno(1,2,3-cd)pyrene	17.00	276	43017	5.314	ng	92
88) Benzo(b)fluoranthene	15.06	252	95151m	12.708	ng	
89) Benzo(k)fluoranthene	15.09	252	28684m	4.000	ng	
90) Benzo(a)pyrene	15.43	252	61234	9.089	ng	# 97
92) Benzo(a,h,i)perylene	17.46	276	39003	7.614	ng	96

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

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