

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF061119\
 Data File : BF114947.D
 Acq On : 11 Jun 2019 15:38
 Operator : HP/JU
 Sample : K3158-01 2X
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 OK-03-060619

Manual Integrations
 APPROVED

mohammad
 6/12/2019 9:03:58 AM

Quant Time: Jun 11 16:42:17 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 10 16:44:00 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.66	152	234540	20.00	ng	0.00
21) Naphthalene-d8	7.95	136	919756	20.00	ng	0.00
39) Acenaphthene-d10	9.69	164	456656	20.00	ng	0.00
64) Phenanthrene-d10	11.16	188	815868	20.00	ng	0.00
76) Chrysene-d12	13.79	240	603623	20.00	ng	0.00
87) Perylene-d12	15.19	264	412046	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.26	112	638028	46.04	ng	0.00
7) Phenol-d6	6.30	99	801957	47.98	ng	0.00
23) Nitrobenzene-d5	7.22	82	458832	30.65	ng	0.00
42) 2,4,6-Tribromophenol	10.47	330	208456	41.96	ng	0.00
45) 2-Fluorobiphenyl	9.02	172	965610	33.56	ng	0.00
79) Terphenyl-d14	12.74	244	962922	31.72	ng	0.00
Target Compounds						
49) Acenaphthylene	9.55	152	121097	2.849	ng	96
50) Dimethylphthalate	9.41	163	90792	2.645	ng	99
71) Phenanthrene	11.19	178	297240	7.509	ng	97
72) Anthracene	11.24	178	98736	2.458	ng	96
75) Fluoranthene	12.37	202	540979	13.213	ng	96
78) Pyrene	12.60	202	584241	11.114	ng	95
81) Benzo(a)anthracene	13.78	228	322907m	7.235	ng	
83) Chrysene	13.82	228	293643	6.552	ng	96
86) Indeno(1,2,3-cd)pyrene	16.50	276	121216	2.877	ng	96
88) Benzo(b)fluoranthene	14.80	252	308103m	11.485	ng	
89) Benzo(k)fluoranthene	14.82	252	105263m	4.522	ng	
90) Benzo(a)pyrene	15.13	252	205418	8.816	ng	# 91
92) Benzo(a,h,i)perylene	16.89	276	124975	6.370	ng	92

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

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