

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF071519\
 Data File : BF115574.D
 Acq On : 15 Jul 2019 23:13
 Operator : HP/JU
 Sample : K3620-01 2X
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CL-01-071219-A

Quant Time: Jul 16 05:20:48 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF071219.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jul 12 14:03:52 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.88	152	205153	20.00	ng	0.00
21) Naphthalene-d8	8.16	136	752742	20.00	ng	0.00
39) Acenaphthene-d10	9.92	164	353843	20.00	ng	-0.01
64) Phenanthrene-d10	11.40	188	583917	20.00	ng	-0.01
76) Chrysene-d12	14.03	240	363986	20.00	ng	-0.02
87) Perylene-d12	15.50	264	401697	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.50	112	602537	48.04	ng	0.00
7) Phenol-d6	6.50	99	789535	49.61	ng	-0.01
23) Nitrobenzene-d5	7.43	82	471013	36.38	ng	-0.02
42) 2,4,6-Tribromophenol	10.70	330	187707	45.45	ng	-0.01
45) 2-Fluorobiphenyl	9.24	172	844845	41.79	ng	-0.01
79) Terphenyl-d14	12.98	244	613734	31.11	ng	-0.02
Target Compounds						
50) Dimethylphthalate	9.63	163	98893	3.715	ng	Qvalue 99
71) Phenanthrene	11.42	178	186308	6.225	ng	96
75) Fluoranthene	12.61	202	182163	5.759	ng	99
78) Pyrene	12.84	202	221088	7.169	ng	97
81) Benzo(a)anthracene	14.02	228	89634	3.738	ng	98
83) Chrysene	14.06	228	95245	3.957	ng	96
88) Benzo(b)fluoranthene	15.07	252	84016m	3.248	ng	
90) Benzo(a)pyrene	15.44	252	70130	3.155	ng	# 95
92) Benzo(g,h,i)perylene	17.42	276	39970	2.053	ng	# 93

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

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