

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF060719\
 Data File : BF114918.D
 Acq On : 8 Jun 2019 4:04
 Operator : HP/JU
 Sample : K3160-05 2X
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 EO-02-060619-C

Quant Time: Jun 08 06:55:35 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF060419.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 04 16:57:15 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.66	152	292570	20.00	ng	0.00
21) Naphthalene-d8	7.95	136	1084674	20.00	ng	0.00
39) Acenaphthene-d10	9.69	164	465289	20.00	ng	-0.01
64) Phenanthrene-d10	11.16	188	757301	20.00	ng	0.00
76) Chrysene-d12	13.78	240	552522	20.00	ng	-0.01
87) Perylene-d12	15.16	264	430603	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.27	112	840994	50.23	ng	0.00
7) Phenol-d6	6.30	99	1034622	50.78	ng	0.00
23) Nitrobenzene-d5	7.22	82	581599	33.01	ng	-0.01
42) 2,4,6-Tribromophenol	10.47	330	205197	40.92	ng	-0.01
45) 2-Fluorobiphenyl	9.02	172	1115988	44.93	ng	-0.01
79) Terphenyl-d14	12.74	244	936822	32.97	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.41	163	97673	2.874	ng	Qvalue 99
71) Phenanthrene	11.19	178	244719	6.769	ng	95
75) Fluoranthene	12.37	202	299606	7.962	ng	90
78) Pyrene	12.59	202	332500	6.937	ng	95
81) Benzo(a)anthracene	13.77	228	145388	3.415	ng	97
83) Chrysene	13.81	228	138549	3.345	ng	95
88) Benzo(b)fluoranthene	14.78	252	119452m	4.358	ng	
90) Benzo(a)pyrene	15.10	252	89921	3.790	ng	# 83
92) Benzo(g,h,i)perylene	16.86	276	51750	2.290	ng	# 92

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

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