

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF061419\
 Data File : BF115057.D
 Acq On : 14 Jun 2019 23:03
 Operator : HP/JU
 Sample : K3351-05
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 RBR-200029

Quant Time: Jun 15 06:38:47 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061219.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jun 12 15:42:37 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.65	152	312258	20.00	ng	0.00
21) Naphthalene-d8	7.93	136	1186123	20.00	ng	0.00
39) Acenaphthene-d10	9.67	164	522197	20.00	ng	0.00
64) Phenanthrene-d10	11.15	188	864417	20.00	ng	0.00
76) Chrysene-d12	13.77	240	535726	20.00	ng	-0.01
87) Perylene-d12	15.14	264	591898	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.26	112	1495789	79.99	ng	0.00
7) Phenol-d6	6.29	99	1827196	81.00	ng	0.00
23) Nitrobenzene-d5	7.21	82	1091639	55.40	ng	0.00
42) 2,4,6-Tribromophenol	10.46	330	390536	69.84	ng	-0.01
45) 2-Fluorobiphenyl	9.00	172	1784744	57.93	ng	0.00
79) Terphenyl-d14	12.73	244	1373073	48.21	ng	-0.01
Target Compounds				Qvalue		
50) Dimethylphthalate	9.40	163	221734	5.646	ng	98
52) Acenaphthene	9.70	154	386694	12.991	ng	97
55) Dibenzofuran	9.87	168	90777	2.023	ng	96
71) Phenanthrene	11.17	178	881524	19.751	ng	98
72) Anthracene	11.22	178	231280	5.137	ng	98
75) Fluoranthene	12.36	202	2557828	56.047	ng	100
78) Pyrene	12.59	202	1939891	39.749	ng	99
81) Benzo(a)anthracene	13.76	228	617175	15.395	ng	98
83) Chrysene	13.80	228	636308	15.961	ng	96
86) Indeno(1,2,3-cd)pyrene	16.44	276	100109	2.859	ng	96
88) Benzo(b)fluoranthene	14.77	252	630291m	16.019	ng	
89) Benzo(k)fluoranthene	14.79	252	213025m	6.168	ng	
90) Benzo(a)pyrene	15.09	252	259239	7.635	ng	99
92) Benzo(a,h,i)perylene	16.83	276	78811	2.643	ng	96

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

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