

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF091019\
 Data File : BF116940.D
 Acq On : 10 Sep 2019 23:48
 Operator : HP/JU
 Sample : K4689-02DL 5X
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MW-2ADL

Quant Time: Sep 11 03:58:39 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF083019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Sep 10 17:49:40 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.83	152	73768	20.00	ng	0.00
21) Naphthalene-d8	8.12	136	296331	20.00	ng	0.00
39) Acenaphthene-d10	9.87	164	157983	20.00	ng	0.00
64) Phenanthrene-d10	11.35	188	243431	20.00	ng	0.00
76) Chrysene-d12	13.98	240	152301	20.00	ng	0.00
87) Perylene-d12	15.43	264	169947	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.45	112	53521	11.75	ng	0.00
7) Phenol-d6	6.46	99	46152	7.72	ng	-0.01
23) Nitrobenzene-d5	7.39	82	96649	18.62	ng	0.00
42) 2,4,6-Tribromophenol	10.65	330	31593	29.91	ng	0.00
45) 2-Fluorobiphenyl	9.19	172	185698	19.83	ng	0.00
79) Terphenyl-d14	12.93	244	152933	18.58	ng	0.00
Target Compounds						Qvalue
27) 2,4-Dimethylphenol	7.77	122	17566	4.431	ng	95
31) Naphthalene	8.16	128	4619686	327.850	ng	95
37) 2-Methylnaphthalene	8.83	142	1132593	120.805	ng	98
38) 1-Methylnaphthalene	8.94	142	1434008	162.029	ng	98
46) 1,1'-Biphenyl	9.29	154	270720	22.583	ng	97
52) Acenaphthene	9.90	154	412973	46.829	ng	99
55) Dibenzofuran	10.07	168	389028	31.292	ng	97
58) Fluorene	10.42	166	220300	23.256	ng	99
71) Phenanthrene	11.37	178	240414	17.742	ng	100
72) Anthracene	11.42	178	36609	2.684	ng	99
73) Carbazole	11.58	167	63970	4.923	ng	98
75) Fluoranthene	12.56	202	27353	2.054	ng	97

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

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