

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF100820\
 Data File : BF121887.D
 Acq On : 8 Oct 2020 14:08
 Operator : JU/CG
 Sample : L4309-01
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 OWLS-HEAD-BH-02-A

Quant Time: Oct 08 14:55:30 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF091020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 01 14:29:14 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.76	152	112527	20.00	ng	-0.01
21) Naphthalene-d8	8.04	136	453441	20.00	ng	-0.02
39) Acenaphthene-d10	9.79	164	238657	20.00	ng	-0.02
64) Phenanthrene-d10	11.28	188	445095	20.00	ng	-0.01
76) Chrysene-d12	13.92	240	327434	20.00	ng	-0.01
86) Perylene-d12	15.36	264	291142	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.38	112	551464	72.83	ng	0.00
7) Phenol-d6	6.40	99	715191	72.00	ng	-0.01
23) Nitrobenzene-d5	7.32	82	491235	58.17	ng	-0.02
42) 2,4,6-Tribromophenol	10.59	330	218938	73.81	ng	-0.02
45) 2-Fluorobiphenyl	9.12	172	889693	56.46	ng	-0.02
79) Terphenyl-d14	12.87	244	953803	59.01	ng	-0.01
Target Compounds						Qvalue
50) Dimethylphthalate	9.52	163	104897	6.064	ng	99
75) Fluoranthene	12.49	202	66765	2.579	ng	99
78) Pyrene	12.72	202	62919	2.630	ng	96
84) Bis(2-ethylhexyl)phthalate	13.90	149	30068	2.326	ng	97
87) Indeno(1,2,3-cd)pyrene	16.78	276	45921	2.422	ng	97
88) Benzo(b)fluoranthene	14.95	252	46008	2.364	ng	# 95
90) Benzo(a)pyrene	15.30	252	36405	2.200	ng	# 97
92) Benzo(g,h,i)perylene	17.21	276	45322	2.921	ng	98

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

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