

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF092220\
 Data File : BF121665.D
 Acq On : 23 Sep 2020 4:06
 Operator : JU/CG
 Sample : L4053-05DL 5X
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 1799-1801DL

Quant Time: Sep 23 07:21:51 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF091020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Sep 22 17:53:26 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.81	152	157246	20.00	ng	0.00
21) Naphthalene-d8	8.10	136	536747	20.00	ng	0.00
39) Acenaphthene-d10	9.85	164	294836	20.00	ng	0.00
64) Phenanthrene-d10	11.34	188	447041	20.00	ng	0.00
76) Chrysene-d12	13.97	240	317411	20.00	ng	0.00
86) Perylene-d12	15.43	264	340351	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.43	112	129122	12.20	ng	0.00
7) Phenol-d6	6.45	99	101574	7.32	ng	0.00
23) Nitrobenzene-d5	7.38	82	222278	22.23	ng	0.00
42) 2,4,6-Tribromophenol	10.64	330	83953	22.91	ng	0.00
45) 2-Fluorobiphenyl	9.17	172	402534	20.68	ng	0.00
79) Terphenyl-d14	12.92	244	281795	17.99	ng	0.00
Target Compounds						
10) Phenol	6.46	94	111945	7.573	ng	# 75
20) 3+4-Methylphenols	7.27	107	6800046	617.450	ng	# 57
27) 2,4-Dimethylphenol	7.87	122	193433	21.534	ng	# 5
32) Benzoic acid	8.16	122	3998327	517.317	ng	98

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

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