

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF092120\
 Data File : BF121640.D
 Acq On : 21 Sep 2020 22:12
 Operator : JU/CG
 Sample : L4053-05
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 1799-1801

Manual Integrations
 APPROVED

mohammad
 9/22/2020 10:09:14 AM

Quant Time: Sep 22 05:40:02 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF091020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Sep 21 14:19:01 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.82	152	140510	20.00	ng	0.00
21) Naphthalene-d8	8.12	136	506284	20.00	ng	0.01
39) Acenaphthene-d10	9.87	164	274991	20.00	ng	0.00
64) Phenanthrene-d10	11.36	188	413671	20.00	ng	0.01
76) Chrysene-d12	13.99	240	353148	20.00	ng	0.00
86) Perylene-d12	15.46	264	343170	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.45	112	533666	56.44	ng	0.01
7) Phenol-d6	6.46	99	429904	34.66	ng	0.00
23) Nitrobenzene-d5	7.42	82	932883	98.93	ng	0.04
42) 2,4,6-Tribromophenol	10.67	330	385124	112.68	ng	0.01
45) 2-Fluorobiphenyl	9.19	172	1324735	72.96	ng	0.00
79) Terphenyl-d14	12.95	244	1391149	79.81	ng	0.00
Target Compounds						
10) Phenol	6.48	94	480721	36.394	ng	79
20) 3+4-Methylphenols	7.35	107	17669453	1795.501	ng	# 57
27) 2,4-Dimethylphenol	7.91	122	709531	83.740	ng	# 8
32) Benzoic acid	8.39	122	13727409m	1864.655	ng	
50) Dimethylphthalate	9.60	163	170652	8.562	ng	96

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

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