

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF060721\
 Data File : BF124162.D
 Acq On : 7 Jun 2021 18:02
 Operator : JU/CG
 Sample : M2583-20DL 5X
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 35-38DL

Quant Time: Jun 07 18:23:50 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF060321.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 07 13:32:21 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.857	152	33388	20.000	ng	0.00
21) Naphthalene-d8	8.145	136	135599	20.000	ng	# 0.00
39) Acenaphthene-d10	9.904	164	76138	20.000	ng	0.00
64) Phenanthrene-d10	11.386	188	136369	20.000	ng	0.00
76) Chrysene-d12	14.027	240	85860	20.000	ng	0.00
86) Perylene-d12	15.510	264	88462	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.487	112	27920	12.588	ng	0.00
7) Phenol-d6	6.504	99	23629	7.926	ng	0.00
23) Nitrobenzene-d5	7.416	82	48357	14.144	ng	-0.01
42) 2,4,6-Tribromophenol	10.692	330	20471	18.978	ng	0.00
45) 2-Fluorobiphenyl	9.222	172	87000	14.379	ng	0.00
79) Terphenyl-d14	12.969	244	68874	11.013	ng	0.00
Target Compounds						
						Qvalue
10) Phenol	6.516	94	15390	4.777	ng	83
20) 3+4-Methylphenols	7.281	107	100425	37.778	ng	# 64
27) 2,4-Dimethylphenol	7.828	122	8319	3.851	ng	84
36) 4-Chloro-3-methylphenol	8.710	107	20907	7.644	ng	# 93

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

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