

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF060721\
 Data File : BF124160.D
 Acq On : 7 Jun 2021 16:53
 Operator : JU/CG
 Sample : M2583-19DL 5X
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 33-34DL

Quant Time: Jun 07 17:15:48 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	32711	20.000	ng	0.00
21) Naphthalene-d8	8.145	136	121958	20.000	ng	0.00
39) Acenaphthene-d10	9.898	164	75116	20.000	ng	0.00
64) Phenanthrene-d10	11.386	188	128271	20.000	ng	0.00
76) Chrysene-d12	14.027	240	87826	20.000	ng	0.00
86) Perylene-d12	15.509	264	82198	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.492	112	31692	14.584	ng	0.00
7) Phenol-d6	6.510	99	25424	8.704	ng	0.00
23) Nitrobenzene-d5	7.422	82	58203	18.928	ng	0.00
42) 2,4,6-Tribromophenol	10.692	330	25034	23.524	ng	0.00
45) 2-Fluorobiphenyl	9.222	172	103296	17.304	ng	0.00
79) Terphenyl-d14	12.968	244	81886	12.801	ng	0.00
Target Compounds						
						Qvalue
10) Phenol	6.522	94	21817	6.912	ng	90
20) 3+4-Methylphenols	7.281	107	76773	29.478	ng	# 61
27) 2,4-Dimethylphenol	7.804	122	5539	2.851	ng	# 38
36) 4-Chloro-3-methylphenol	8.710	107	35980	14.626	ng	# 92

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

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