

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM012021\
 Data File : BM028470.D
 Acq On : 20 Jan 2021 16:30
 Operator : CG/JU
 Sample : M1125-07
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 1693-94

Quant Time: Jan 20 17:01:23 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM012021.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jan 20 14:45:01 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.028	152	37317	20.00	ng	0.00
21) Naphthalene-d8	10.851	136	145745	20.00	ng	# 0.00
39) Acenaphthene-d10	14.669	164	79508	20.00	ng	0.00
64) Phenanthrene-d10	17.398	188	152812	20.00	ng	# 0.00
76) Chrysene-d12	21.527	240	140504	20.00	ng	# 0.00
86) Perylene-d12	23.815	264	139134	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.557	112	193772	82.03	ng	0.00
7) Phenol-d6	7.187	99	154702	48.10	ng	0.00
23) Nitrobenzene-d5	9.204	82	247552	80.56	ng	0.00
42) 2,4,6-Tribromophenol	16.151	330	132842	126.10	ng	0.00
45) 2-Fluorobiphenyl	13.298	172	514115	81.13	ng	0.00
79) Terphenyl-d14	19.992	244	587870	75.55	ng	0.00
Target Compounds						
10) Phenol	7.210	94	15353	5.06	ng	Qvalue 93
15) Benzyl Alcohol	8.275	79	12867	5.74	ng	86
32) Benzoic acid	10.181	122	45002	25.62	ng	# 88
50) Dimethylphthalate	14.121	163	23333	3.57	ng	98

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

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