

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM020323\
 Data File : BM038692.D
 Acq On : 03 Feb 2023 23:02
 Operator : CG/JU
 Sample : 01375-01RE
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 MANHOLERE

Quant Time: Feb 04 05:23:43 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM020123.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Feb 04 05:17:53 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.922	152	78255	20.000	ng	0.00
21) Naphthalene-d8	10.733	136	280694	20.000	ng	0.00
39) Acenaphthene-d10	14.563	164	204963	20.000	ng	0.00
64) Phenanthrene-d10	17.304	188	427016	20.000	ng	0.00
76) Chrysene-d12	21.486	240	348055	20.000	ng	0.00
86) Perylene-d12	23.880	264	413999	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.487	112	66406	16.623	ng	0.00
7) Phenol-d6	7.104	99	50356	11.052	ng	0.00
23) Nitrobenzene-d5	9.098	82	241123	48.793	ng	0.00
42) 2,4,6-Tribromophenol	16.051	330	122850	44.097	ng	0.00
45) 2-Fluorobiphenyl	13.186	172	837853	52.042	ng	0.00
79) Terphenyl-d14	19.927	244	1075296	54.519	ng	0.00
Target Compounds						
						Qvalue
10) Phenol	7.128	94	35920	7.804	ng	96
15) Benzyl Alcohol	8.181	79	8022	2.392	ng	91
20) 3+4-Methylphenols	8.704	107	59292	12.596	ng	92
32) Benzoic acid	10.098	122	81315	24.945	ng	97
84) Bis(2-ethylhexyl)phtha...	21.380	149	20280	2.508	ng	# 96

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

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