

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM020323\
 Data File : BM038683.D
 Acq On : 03 Feb 2023 17:34
 Operator : CG/JU
 Sample : 01375-01
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 MANHOLE

Manual Integrations
 APPROVED

Reviewed By :Christian Giraldo 02/06/2023
 Supervised By :Jagrut Upadhyay 02/06/2023

Quant Time: Feb 04 05:21:15 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	74523	20.000	ng	0.00
21) Naphthalene-d8	10.733	136	263133	20.000	ng	0.00
39) Acenaphthene-d10	14.563	164	193294	20.000	ng	0.00
64) Phenanthrene-d10	17.309	188	403010	20.000	ng	0.00
76) Chrysene-d12	21.486	240	323927	20.000	ng	0.00
86) Perylene-d12	23.886	264	389917	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.493	112	63185	16.609	ng	0.01
7) Phenol-d6	7.110	99	47760	11.007	ng	0.01
23) Nitrobenzene-d5	9.098	82	228902	49.411	ng	0.00
42) 2,4,6-Tribromophenol	16.057	330	117402	44.686	ng	0.00
45) 2-Fluorobiphenyl	13.186	172	803364	52.912	ng	0.00
79) Terphenyl-d14	19.927	244	1018928	55.509	ng	0.00
Target Compounds						Qvalue
10) Phenol	7.134	94	33553	7.655	ng	97
15) Benzyl Alcohol	8.181	79	10462m	3.276	ng	
20) 3+4-Methylphenols	8.710	107	55039	12.278	ng	92
32) Benzoic acid	10.098	122	77386	25.324	ng	95
84) Bis(2-ethylhexyl)phtha...	21.386	149	21390	2.643	ng	# 98

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

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