

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM111022\
 Data File : BM037453.D
 Acq On : 10 Nov 2022 16:09
 Operator : CG/JU
 Sample : N5436-09DL 20X
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 250FERRYBLVD-TFS-09DL

Quant Time: Nov 10 16:56:45 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM102822.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 09 02:08:14 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.810	152	248196	20.000	ng	-0.01
21) Naphthalene-d8	10.610	136	959222	20.000	ng	-0.02
39) Acenaphthene-d10	14.451	164	551630	20.000	ng	-0.01
64) Phenanthrene-d10	17.192	188	1123827	20.000	ng	-0.01
76) Chrysene-d12	21.374	240	1236120	20.000	ng	0.00
86) Perylene-d12	23.703	264	1325247	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.387	112	78943	5.495	ng	0.00
7) Phenol-d6	6.992	99	93817	4.812	ng	-0.01
23) Nitrobenzene-d5	8.980	82	58281	3.065	ng	-0.02
42) 2,4,6-Tribromophenol	15.939	330	36438	5.605	ng	-0.02
45) 2-Fluorobiphenyl	13.074	172	132005	3.180	ng	-0.02
79) Terphenyl-d14	19.815	244	223955	3.249	ng	-0.01
Target Compounds						
						Qvalue
20) 3+4-Methylphenols	8.592	107	346775	17.682	ng	92
27) 2,4-Dimethylphenol	9.798	122	561688	43.860	ng	99
75) Fluoranthene	19.250	202	175997	2.388	ng	98
78) Pyrene	19.609	202	292106	3.073	ng	# 91

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

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