

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM052522\
 Data File : BM035253.D
 Acq On : 25 May 2022 22:10
 Operator : CG/JU
 Sample : N2895-06
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 P030-SS001-0612-01

Quant Time: May 26 05:46:24 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM042722.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu May 26 05:05:10 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.886	152	113524	20.000	ng	0.00
21) Naphthalene-d8	10.686	136	430129	20.000	ng	0.00
39) Acenaphthene-d10	14.521	164	273306	20.000	ng	0.00
64) Phenanthrene-d10	17.268	188	612575	20.000	ng	# 0.00
76) Chrysene-d12	21.450	240	656234	20.000	ng	# 0.00
86) Perylene-d12	23.827	264	705400	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.469	112	579206	90.012	ng	0.00
7) Phenol-d6	7.057	99	736233	88.629	ng	0.00
23) Nitrobenzene-d5	9.045	82	463053	58.135	ng	0.00
42) 2,4,6-Tribromophenol	16.015	330	328166	110.381	ng	0.00
45) 2-Fluorobiphenyl	13.151	172	1107203	58.809	ng	0.00
79) Terphenyl-d14	19.892	244	1911635	56.402	ng	0.00

Target Compounds	Qvalue

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

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