

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM052522\
 Data File : BM035254.D
 Acq On : 25 May 2022 22:46
 Operator : CG/JU
 Sample : N2895-07
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 P030-SS001-1218-01

Quant Time: May 26 05:48:20 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	7.887	152	113947	20.000	ng	0.00
21) Naphthalene-d8	10.686	136	423496	20.000	ng	0.00
39) Acenaphthene-d10	14.521	164	269541	20.000	ng	0.00
64) Phenanthrene-d10	17.268	188	583324	20.000	ng	0.00
76) Chrysene-d12	21.450	240	622587	20.000	ng	# 0.00
86) Perylene-d12	23.827	264	681073	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.469	112	627015	97.080	ng	0.00
7) Phenol-d6	7.057	99	788205	94.533	ng	0.00
23) Nitrobenzene-d5	9.045	82	504341	64.311	ng	0.00
42) 2,4,6-Tribromophenol	16.015	330	336005	114.597	ng	0.00
45) 2-Fluorobiphenyl	13.151	172	1239793	66.772	ng	0.00
79) Terphenyl-d14	19.892	244	2060779	64.089	ng	0.00

Target Compounds	Qvalue

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

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