

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM081023\
 Data File : BM041369.D
 Acq On : 10 Aug 2023 19:02
 Operator : MA/JU
 Sample : 03955-17
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 WC-27

Quant Time: Aug 11 03:35:17 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM072823.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 09 12:20:12 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.775	152	168194	20.000	ng	0.00
21) Naphthalene-d8	10.580	136	644583	20.000	ng	0.00
39) Acenaphthene-d10	14.439	164	396647	20.000	ng	0.00
64) Phenanthrene-d10	17.203	188	812634	20.000	ng	# 0.00
76) Chrysene-d12	21.415	240	603209	20.000	ng	0.00
86) Perylene-d12	23.780	264	697800	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.351	112	1305916	116.045	ng	0.00
7) Phenol-d6	6.963	99	1484786	101.786	ng	0.00
23) Nitrobenzene-d5	8.951	82	1136437	81.990	ng	0.00
42) 2,4,6-Tribromophenol	15.945	330	738783	131.682	ng	0.00
45) 2-Fluorobiphenyl	13.057	172	2624785	83.595	ng	0.00
79) Terphenyl-d14	19.845	244	3570849	107.032	ng	0.00

Target Compounds Qvalue

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

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