

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM111023\
 Data File : BM042689.D
 Acq On : 10 Nov 2023 13:46
 Operator : MA/JU
 Sample : 05295-04
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 TP-2

Quant Time: Nov 10 14:29:27 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM103023.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Nov 09 13:21:00 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.845	152	36774	20.000	ng	0.00
21) Naphthalene-d8	10.663	136	140638	20.000	ng	0.00
39) Acenaphthene-d10	14.504	164	102302	20.000	ng	0.00
64) Phenanthrene-d10	17.256	188	243029	20.000	ng	# 0.00
76) Chrysene-d12	21.439	240	265780	20.000	ng	# 0.00
86) Perylene-d12	23.815	264	290266	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.381	112	367778	161.405	ng	0.00
7) Phenol-d6	7.016	99	447769	143.285	ng	0.00
23) Nitrobenzene-d5	9.051	82	386457	128.097	ng	0.00
42) 2,4,6-Tribromophenol	15.998	330	324696	243.101	ng	0.00
45) 2-Fluorobiphenyl	13.121	172	900107	118.393	ng	0.00
79) Terphenyl-d14	19.874	244	2214566	141.284	ng	0.00

Target Compounds	Qvalue

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

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