

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110821\
 Data File : BM032891.D
 Acq On : 08 Nov 2021 12:51
 Operator : CG/JU
 Sample : M4499-02
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 WEL-BRIDGE

Quant Time: Nov 08 13:25:46 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM110221.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 02 16:07:18 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.490	152	85418	20.000	ng	0.00
21) Naphthalene-d8	10.266	136	357996	20.000	ng	# 0.00
39) Acenaphthene-d10	14.148	164	250556	20.000	ng	0.00
64) Phenanthrene-d10	16.883	188	573680	20.000	ng	-0.01
76) Chrysene-d12	21.065	240	617991	20.000	ng	-0.01
86) Perylene-d12	23.177	264	633288	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.084	112	789010	161.038	ng	0.00
7) Phenol-d6	6.696	99	986519	135.207	ng	0.00
23) Nitrobenzene-d5	8.642	82	655323	100.190	ng	0.00
42) 2,4,6-Tribromophenol	15.642	330	434873	154.976	ng	0.00
45) 2-Fluorobiphenyl	12.760	172	1556140	96.328	ng	0.00
79) Terphenyl-d14	19.518	244	2588858	91.460	ng	0.00

Target Compounds	Qvalue

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

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