

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110821\
 Data File : BM032899.D
 Acq On : 08 Nov 2021 17:41
 Operator : CG/JU
 Sample : M4475-06
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DUP01

Quant Time: Nov 09 00:24:38 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	7.484	152	85261	20.000	ng	0.00
21) Naphthalene-d8	10.266	136	375289	20.000	ng	# 0.00
39) Acenaphthene-d10	14.148	164	270915	20.000	ng	0.00
64) Phenanthrene-d10	16.883	188	600026	20.000	ng	-0.01
76) Chrysene-d12	21.065	240	637911	20.000	ng	-0.01
86) Perylene-d12	23.171	264	656589	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.078	112	575955	117.770	ng	0.00
7) Phenol-d6	6.689	99	767054	105.322	ng	0.00
23) Nitrobenzene-d5	8.636	82	458169	66.820	ng	-0.01
42) 2,4,6-Tribromophenol	15.642	330	310216	102.244	ng	0.00
45) 2-Fluorobiphenyl	12.760	172	1111444	63.630	ng	0.00
79) Terphenyl-d14	19.512	244	1727507	59.124	ng	-0.01

Target Compounds	Qvalue

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

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