

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110421\
 Data File : BM032875.D
 Acq On : 04 Nov 2021 15:12
 Operator : CG/JU
 Sample : M4472-07
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 WC-2-KD

Quant Time: Nov 07 08:47:43 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	110658	20.000	ng	0.00
21) Naphthalene-d8	10.266	136	422902	20.000	ng	# 0.00
39) Acenaphthene-d10	14.148	164	245893	20.000	ng	0.00
64) Phenanthrene-d10	16.883	188	473661	20.000	ng	-0.01
76) Chrysene-d12	21.065	240	445942	20.000	ng	-0.01
86) Perylene-d12	23.177	264	477073	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.078	112	579487	91.297	ng	0.00
7) Phenol-d6	6.690	99	719666	76.136	ng	0.00
23) Nitrobenzene-d5	8.637	82	455057	58.894	ng	-0.01
42) 2,4,6-Tribromophenol	15.642	330	225934	82.043	ng	0.00
45) 2-Fluorobiphenyl	12.760	172	955615	60.276	ng	0.00
79) Terphenyl-d14	19.512	244	1195023	58.506	ng	-0.01

Target Compounds	Qvalue

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

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