

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110121\
 Data File : BM032840.D
 Acq On : 01 Nov 2021 19:39
 Operator : CG/JU
 Sample : PB140347BL
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB140347BL

Quant Time: Nov 02 00:07:42 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM110121.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Nov 01 18:47:37 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.495	152	129081	20.000	ng	0.00
21) Naphthalene-d8	10.278	136	559454	20.000	ng	# 0.00
39) Acenaphthene-d10	14.154	164	374028	20.000	ng	0.00
64) Phenanthrene-d10	16.889	188	788702	20.000	ng	0.00
76) Chrysene-d12	21.071	240	741749	20.000	ng	#-0.01
86) Perylene-d12	23.188	264	728357	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.090	112	1040737	147.640	ng	0.00
7) Phenol-d6	6.701	99	1469167	132.742	ng	0.00
23) Nitrobenzene-d5	8.648	82	950834	87.359	ng	0.00
42) 2,4,6-Tribromophenol	15.648	330	494188	122.085	ng	0.00
45) 2-Fluorobiphenyl	12.771	172	2112316	86.245	ng	0.00
79) Terphenyl-d14	19.524	244	2893282	80.555	ng	0.00

Target Compounds	Qvalue

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

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