

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM090921\
 Data File : BM032109.D
 Acq On : 09 Sep 2021 13:14
 Operator : CG/JU
 Sample : M3644-02
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 191-0902-W

Quant Time: Sep 09 14:44:03 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM090821.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Sep 08 19:23:17 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.357	152	63987	20.000	ng	0.00
21) Naphthalene-d8	10.110	136	262451	20.000	ng	# 0.00
39) Acenaphthene-d10	14.015	164	173178	20.000	ng	0.00
64) Phenanthrene-d10	16.780	188	339787	20.000	ng	0.00
76) Chrysene-d12	20.997	240	311251	20.000	ng	-0.01
86) Perylene-d12	23.091	264	319264	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.004	112	147389	41.102	ng	0.00
7) Phenol-d6	6.563	99	118586	21.738	ng	0.00
23) Nitrobenzene-d5	8.510	82	540153	91.722	ng	0.00
42) 2,4,6-Tribromophenol	15.527	330	118648	86.209	ng	0.00
45) 2-Fluorobiphenyl	12.621	172	1134301	85.880	ng	0.00
79) Terphenyl-d14	19.445	244	1635691	85.744	ng	0.00

Target Compounds	Qvalue

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

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