

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110121\
 Data File : BP007795.D
 Acq On : 01 Nov 2021 17:15
 Operator : CG/JU
 Sample : M4424-08
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 T3-EP01-SE

Quant Time: Nov 02 01:49:51 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP101321.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 02 01:43:06 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.928	152	153490	20.000	ng	0.00
21) Naphthalene-d8	10.728	136	582046	20.000	ng	0.00
39) Acenaphthene-d10	14.563	164	381218	20.000	ng	-0.01
64) Phenanthrene-d10	17.322	188	903399	20.000	ng	0.00
76) Chrysene-d12	21.410	240	1123070	20.000	ng	-0.01
86) Perylene-d12	23.851	264	1252114	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.505	112	924949	99.879	ng	0.00
7) Phenol-d6	7.093	99	1150284	92.194	ng	0.00
23) Nitrobenzene-d5	9.081	82	688304	72.980	ng	-0.01
42) 2,4,6-Tribromophenol	16.057	330	537508	135.838	ng	0.00
45) 2-Fluorobiphenyl	13.193	172	1681364	67.415	ng	0.00
79) Terphenyl-d14	19.880	244	3304955	59.603	ng	-0.01

Target Compounds	Qvalue

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

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