

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110121\
 Data File : BP007797.D
 Acq On : 01 Nov 2021 18:23
 Operator : CG/JU
 Sample : M4424-07
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 T3-EP03-SW

Quant Time: Nov 02 01:50:22 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	155682	20.000	ng	0.00
21) Naphthalene-d8	10.728	136	568113	20.000	ng	0.00
39) Acenaphthene-d10	14.563	164	357657	20.000	ng	-0.01
64) Phenanthrene-d10	17.316	188	806587	20.000	ng	-0.01
76) Chrysene-d12	21.410	240	1017914	20.000	ng	-0.01
86) Perylene-d12	23.851	264	1136179	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.499	112	852374	90.747	ng	0.00
7) Phenol-d6	7.093	99	1034096	81.715	ng	0.00
23) Nitrobenzene-d5	9.081	82	622108	67.579	ng	-0.01
42) 2,4,6-Tribromophenol	16.057	330	431086	116.120	ng	0.00
45) 2-Fluorobiphenyl	13.187	172	1465341	62.624	ng	-0.01
79) Terphenyl-d14	19.880	244	2610807	51.949	ng	-0.01

Target Compounds	Qvalue

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

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