

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110421\
 Data File : BP007854.D
 Acq On : 04 Nov 2021 18:36
 Operator : CG/JU
 Sample : M4497-08
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 346

Quant Time: Nov 07 11:17:30 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP110221.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 02 19:18:55 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.910	152	178764	20.000	ng	-0.01
21) Naphthalene-d8	10.710	136	676778	20.000	ng	-0.02
39) Acenaphthene-d10	14.551	164	436130	20.000	ng	-0.01
64) Phenanthrene-d10	17.304	188	936509	20.000	ng	-0.02
76) Chrysene-d12	21.398	240	1011545	20.000	ng	-0.02
86) Perylene-d12	23.833	264	1102552	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.493	112	1517699	143.984	ng	0.00
7) Phenol-d6	7.087	99	1703766	111.674	ng	0.00
23) Nitrobenzene-d5	9.069	82	1182868	93.887	ng	-0.02
42) 2,4,6-Tribromophenol	16.045	330	893011	143.474	ng	-0.01
45) 2-Fluorobiphenyl	13.175	172	2795517	93.345	ng	-0.02
79) Terphenyl-d14	19.875	244	4751687	95.654	ng	0.00

Target Compounds	Qvalue

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

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