

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110821\
 Data File : BP007877.D
 Acq On : 09 Nov 2021 00:20
 Operator : CG/JU
 Sample : M4477-10
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 T5-B1-NE

Quant Time: Nov 09 08:47:54 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP110221.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Nov 08 16:36:40 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.863	152	123498	20.000	ng	0.00
21) Naphthalene-d8	10.657	136	432089	20.000	ng	-0.01
39) Acenaphthene-d10	14.498	164	287602	20.000	ng	0.00
64) Phenanthrene-d10	17.251	188	645507	20.000	ng	-0.01
76) Chrysene-d12	21.339	240	799095	20.000	ng	-0.01
86) Perylene-d12	23.750	264	899037	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.446	112	712446	97.836	ng	0.00
7) Phenol-d6	7.034	99	895692	84.981	ng	-0.01
23) Nitrobenzene-d5	9.016	82	499724	62.126	ng	-0.02
42) 2,4,6-Tribromophenol	15.986	330	417351	101.681	ng	-0.01
45) 2-Fluorobiphenyl	13.122	172	1261614	63.882	ng	-0.01
79) Terphenyl-d14	19.815	244	2383431	60.736	ng	0.00
Target Compounds						
						Qvalue
71) Phenanthrene	17.292	178	148404	4.373	ng	99
75) Fluoranthene	19.263	202	165518	3.867	ng	99
78) Pyrene	19.615	202	165460	3.376	ng	99

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

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