

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110821\
 Data File : BP007879.D
 Acq On : 09 Nov 2021 01:28
 Operator : CG/JU
 Sample : M4477-09
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 T5-EP3-N

Quant Time: Nov 09 08:48:41 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	155790	20.000	ng	0.00
21) Naphthalene-d8	10.657	136	540027	20.000	ng	-0.01
39) Acenaphthene-d10	14.498	164	358173	20.000	ng	0.00
64) Phenanthrene-d10	17.251	188	798236	20.000	ng	#-0.01
76) Chrysene-d12	21.339	240	929910	20.000	ng	-0.01
86) Perylene-d12	23.751	264	1035679	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.446	112	1018725	110.899	ng	0.00
7) Phenol-d6	7.040	99	1257936	94.611	ng	0.00
23) Nitrobenzene-d5	9.016	82	712812	70.904	ng	-0.02
42) 2,4,6-Tribromophenol	15.986	330	588592	115.147	ng	-0.01
45) 2-Fluorobiphenyl	13.122	172	1758308	71.490	ng	-0.01
79) Terphenyl-d14	19.816	244	3141340	68.788	ng	0.00
Target Compounds						
						Qvalue
75) Fluoranthene	19.263	202	110948	2.096	ng	97
78) Pyrene	19.610	202	115110	2.018	ng	98
88) Benzo(b)fluoranthene	23.015	252	126530	2.073	ng	# 92

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

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