

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP012722\
 Data File : BP008956.D
 Acq On : 27 Jan 2022 20:50
 Operator : CG/JU
 Sample : N1281-01
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BRICK-SAMPLE

Quant Time: Jan 28 03:54:48 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP011422.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jan 26 01:52:51 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.828	152	244403	20.000	ng	-0.01
21) Naphthalene-d8	10.628	136	860468	20.000	ng	-0.01
39) Acenaphthene-d10	14.469	164	498802	20.000	ng	-0.01
64) Phenanthrene-d10	17.227	188	1010744	20.000	ng	-0.01
76) Chrysene-d12	21.321	240	1179742	20.000	ng	-0.01
86) Perylene-d12	23.733	264	1354269	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.422	112	785821	49.721	ng	0.00
7) Phenol-d6	7.010	99	1184533	61.893	ng	-0.01
23) Nitrobenzene-d5	8.987	82	821230	54.185	ng	-0.01
42) 2,4,6-Tribromophenol	15.969	330	222001	30.576	ng	-0.01
45) 2-Fluorobiphenyl	13.092	172	2023992	53.511	ng	-0.01
79) Terphenyl-d14	19.786	244	3216013	46.014	ng	-0.02
Target Compounds						
71) Phenanthrene	17.269	178	115282	2.026	ng	98

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

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