

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020122\
 Data File : BP008997.D
 Acq On : 01 Feb 2022 20:56
 Operator : CG/JU
 Sample : N1325-10 5X
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 TES-4

Quant Time: Feb 02 04:07:37 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP011422.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Feb 01 04:16:03 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.828	152	264868	20.000	ng	0.00
21) Naphthalene-d8	10.622	136	1034039	20.000	ng	-0.01
39) Acenaphthene-d10	14.463	164	669738	20.000	ng	-0.01
64) Phenanthrene-d10	17.222	188	1356204	20.000	ng	-0.01
76) Chrysene-d12	21.316	240	1310430	20.000	ng	-0.01
86) Perylene-d12	23.721	264	1333867	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.416	112	240739	14.055	ng	0.00
7) Phenol-d6	7.010	99	304763	14.694	ng	0.00
23) Nitrobenzene-d5	8.987	82	184349	10.122	ng	-0.01
42) 2,4,6-Tribromophenol	15.963	330	140778	14.441	ng	-0.01
45) 2-Fluorobiphenyl	13.087	172	522058	10.280	ng	-0.01
79) Terphenyl-d14	19.780	244	768759	9.902	ng	-0.02
Target Compounds						
						Qvalue
71) Phenanthrene	17.263	178	202771	2.656	ng	97
75) Fluoranthene	19.233	202	316723	3.730	ng	99
78) Pyrene	19.586	202	577398	6.315	ng	96
81) Benzo(a)anthracene	21.298	228	277555	3.148	ng	100
83) Chrysene	21.351	228	310908	3.638	ng	98
88) Benzo(b)fluoranthene	22.986	252	256536	2.880	ng	98
90) Benzo(a)pyrene	23.615	252	208410	2.424	ng	98

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

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