

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP033122\
 Data File : BP009650.D
 Acq On : 31 Mar 2022 21:13
 Operator : CG/JU
 Sample : N2150-01 10X
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 279

Quant Time: Apr 01 04:12:07 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP031722.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Mar 31 04:36:51 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.740	152	336063	20.000	ng	0.00
21) Naphthalene-d8	10.534	136	1356700	20.000	ng	0.00
39) Acenaphthene-d10	14.387	164	867066	20.000	ng	0.00
64) Phenanthrene-d10	17.157	188	1742713	20.000	ng	0.00
76) Chrysene-d12	21.275	240	1361845	20.000	ng	0.00
86) Perylene-d12	23.663	264	1500463	20.000	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.352	112	195840	10.174	ng	0.00
7) Phenol-d6	6.940	99	240836	9.252	ng	0.00
23) Nitrobenzene-d5	8.905	82	143637	5.626	ng	0.00
42) 2,4,6-Tribromophenol	15.898	330	102442	7.160	ng	0.00
45) 2-Fluorobiphenyl	13.010	172	376571	6.021	ng	0.00
79) Terphenyl-d14	19.733	244	475468	6.267	ng	0.00
Target Compounds						
						Qvalue
71) Phenanthrene	17.198	178	403960	4.330	ng	100
75) Fluoranthene	19.186	202	801199	7.094	ng	98
78) Pyrene	19.539	202	876150	9.763	ng	99
81) Benzo(a)anthracene	21.263	228	398964	4.459	ng	96
83) Chrysene	21.310	228	400377	4.726	ng	98
88) Benzo(b)fluoranthene	22.939	252	451178	5.037	ng	# 97
90) Benzo(a)pyrene	23.551	252	265308	3.462	ng	# 95

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

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