

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092322\
 Data File : BP011782.D
 Acq On : 23 Sep 2022 12:55
 Operator : CG/JU
 Sample : N4609-04
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DSN002

Quant Time: Sep 24 01:33:53 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP092122.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Sep 22 02:45:14 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.916	152	38813	20.000	ng	0.00
21) Naphthalene-d8	10.728	136	203690	20.000	ng	0.00
39) Acenaphthene-d10	14.557	164	138038	20.000	ng	-0.01
64) Phenanthrene-d10	17.316	188	373131	20.000	ng	0.00
76) Chrysene-d12	21.398	240	498163	20.000	ng	0.00
86) Perylene-d12	23.851	264	655103	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.475	112	194708	88.021	ng	0.00
7) Phenol-d6	7.075	99	191700	58.610	ng	0.00
23) Nitrobenzene-d5	9.081	82	342704	88.062	ng	-0.01
42) 2,4,6-Tribromophenol	16.051	330	161064	173.285	ng	0.00
45) 2-Fluorobiphenyl	13.181	172	740787	84.866	ng	-0.01
79) Terphenyl-d14	19.869	244	1899815	87.688	ng	0.00
Target Compounds						
10) Phenol	7.099	94	9305	2.526	ng	82
84) Bis(2-ethylhexyl)phtha...	21.310	149	22679	2.812	ng	93

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

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