

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091922\
 Data File : BP011740.D
 Acq On : 19 Sep 2022 18:04
 Operator : CG/JU
 Sample : N4653-02
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 CAR-1948

Quant Time: Sep 20 04:46:09 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP091422.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Sep 20 04:42:52 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.940	152	531994	20.000	ng	0.00
21) Naphthalene-d8	10.752	136	2199958	20.000	ng	# 0.00
39) Acenaphthene-d10	14.575	164	1412589	20.000	ng	0.00
64) Phenanthrene-d10	17.328	188	3153802	20.000	ng	0.00
76) Chrysene-d12	21.404	240	3397348	20.000	ng	0.00
86) Perylene-d12	23.857	264	3432118	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.505	112	4735386	162.020	ng	0.01
7) Phenol-d6	7.105	99	5681604	145.832	ng	0.01
23) Nitrobenzene-d5	9.105	82	3709286	99.996	ng	0.00
42) 2,4,6-Tribromophenol	16.069	330	2838752	257.201	ng	0.00
45) 2-Fluorobiphenyl	13.204	172	8183030	90.048	ng	0.00
79) Terphenyl-d14	19.880	244	13391226	85.715	ng	0.00

Target Compounds	Qvalue

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

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