

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092822\
 Data File : BP011834.D
 Acq On : 28 Sep 2022 17:28
 Operator : CG/JU
 Sample : N4812-02
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 WC-3

Quant Time: Sep 29 02:28:28 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP092122.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Sep 28 02:03:33 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.905	152	392679	20.000	ng	0.00
21) Naphthalene-d8	10.716	136	1596699	20.000	ng	# 0.00
39) Acenaphthene-d10	14.551	164	982448	20.000	ng	0.00
64) Phenanthrene-d10	17.304	188	2134625	20.000	ng	# 0.00
76) Chrysene-d12	21.386	240	2023477	20.000	ng	#-0.01
86) Perylene-d12	23.821	264	1871866	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.475	112	3225457	144.122	ng	0.00
7) Phenol-d6	7.069	99	3820550	115.455	ng	0.00
23) Nitrobenzene-d5	9.075	82	2418680	79.285	ng	0.00
42) 2,4,6-Tribromophenol	16.040	330	1752526	264.921	ng	0.00
45) 2-Fluorobiphenyl	13.175	172	5776530	92.981	ng	0.00
79) Terphenyl-d14	19.857	244	9499249	107.943	ng	-0.01

Target Compounds	Qvalue

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

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