

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP090722\
 Data File : BP011649.D
 Acq On : 07 Sep 2022 15:47
 Operator : CG/JU
 Sample : N4501-12
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SB-13-COMP

Quant Time: Sep 08 04:05:49 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP090222.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Sep 03 04:07:09 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.969	152	748116	20.000	ng	0.00
21) Naphthalene-d8	10.781	136	3078751	20.000	ng	#-0.01
39) Acenaphthene-d10	14.610	164	1752309	20.000	ng	0.00
64) Phenanthrene-d10	17.363	188	4092271	20.000	ng	0.00
76) Chrysene-d12	21.445	240	3336522	20.000	ng	0.00
86) Perylene-d12	23.916	264	2961506	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.528	112	6770902	160.069	ng	0.00
7) Phenol-d6	7.128	99	8566228	153.046	ng	0.00
23) Nitrobenzene-d5	9.134	82	4916682	98.028	ng	0.00
42) 2,4,6-Tribromophenol	16.098	330	3797977	191.833	ng	0.00
45) 2-Fluorobiphenyl	13.240	172	10396083	84.839	ng	0.00
79) Terphenyl-d14	19.916	244	13807180	86.674	ng	0.00

Target Compounds	Qvalue

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

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