

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP090722\
 Data File : BP011648.D
 Acq On : 07 Sep 2022 15:12
 Operator : CG/JU
 Sample : N4501-06
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SB-10-COMP

Quant Time: Sep 08 04:05:22 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	604833	20.000	ng	0.00
21) Naphthalene-d8	10.781	136	2479242	20.000	ng	#-0.01
39) Acenaphthene-d10	14.610	164	1680249	20.000	ng	0.00
64) Phenanthrene-d10	17.363	188	3365852	20.000	ng	0.00
76) Chrysene-d12	21.439	240	2885401	20.000	ng	-0.01
86) Perylene-d12	23.916	264	2598480	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.522	112	4219449	123.381	ng	0.00
7) Phenol-d6	7.122	99	5426652	119.922	ng	0.00
23) Nitrobenzene-d5	9.134	82	3009881	74.522	ng	0.00
42) 2,4,6-Tribromophenol	16.098	330	2430171	128.010	ng	0.00
45) 2-Fluorobiphenyl	13.234	172	7760350	66.046	ng	-0.01
79) Terphenyl-d14	19.916	244	10443061	75.805	ng	0.00

Target Compounds	Qvalue

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

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