

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092324\
 Data File : BP021976.D
 Acq On : 23 Sep 2024 10:29
 Operator : RC/JU
 Sample : PB163138BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB163138BL

Quant Time: Sep 23 11:06:28 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP091924.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Sep 20 01:22:00 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.681	152	102576	20.000	ng	-0.04
21) Naphthalene-d8	10.451	136	359434	20.000	ng	-0.02
39) Acenaphthene-d10	14.316	164	240083	20.000	ng	-0.01
64) Phenanthrene-d10	17.110	188	538548	20.000	ng	-0.03
76) Chrysene-d12	21.533	240	573086	20.000	ng	-0.03
86) Perylene-d12	24.815	264	627189	20.000	ng	-0.05
System Monitoring Compounds						
5) 2-Fluorophenol	5.299	112	767194	140.556	ng	-0.04
7) Phenol-d6	6.863	99	953856	134.615	ng	-0.04
23) Nitrobenzene-d5	8.828	82	637116	93.792	ng	-0.03
42) 2,4,6-Tribromophenol	15.828	330	549886	123.546	ng	-0.03
45) 2-Fluorobiphenyl	12.928	172	1563201	95.047	ng	-0.01
79) Terphenyl-d14	19.833	244	3064736	100.457	ng	-0.02

Target Compounds	Qvalue

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

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