

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP090722\
 Data File : BP011647.D
 Acq On : 07 Sep 2022 14:38
 Operator : CG/JU
 Sample : N4501-04
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SB-08-COMP

Quant Time: Sep 08 04:04:51 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	570449	20.000	ng	0.00
21) Naphthalene-d8	10.781	136	2380546	20.000	ng	#-0.01
39) Acenaphthene-d10	14.604	164	1402030	20.000	ng	-0.01
64) Phenanthrene-d10	17.363	188	2746853	20.000	ng	0.00
76) Chrysene-d12	21.439	240	2340106	20.000	ng	-0.01
86) Perylene-d12	23.910	264	2142906	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.528	112	5889116	182.584	ng	0.00
7) Phenol-d6	7.128	99	7671743	179.754	ng	0.00
23) Nitrobenzene-d5	9.134	82	4361048	112.452	ng	0.00
42) 2,4,6-Tribromophenol	16.098	330	2425191	153.098	ng	0.00
45) 2-Fluorobiphenyl	13.234	172	9626733	98.188	ng	-0.01
79) Terphenyl-d14	19.916	244	12334747	110.401	ng	0.00

Target Compounds Qvalue

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

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