

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP013122\
 Data File : BP008961.D
 Acq On : 31 Jan 2022 12:44
 Operator : CG/JU
 Sample : PB142358BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB142358BL

Quant Time: Feb 01 04:18:48 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP011422.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Feb 01 04:16:03 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.828	152	265787	20.000	ng	0.00
21) Naphthalene-d8	10.628	136	1028478	20.000	ng	0.00
39) Acenaphthene-d10	14.469	164	684091	20.000	ng	0.00
64) Phenanthrene-d10	17.228	188	1432251	20.000	ng	0.00
76) Chrysene-d12	21.316	240	1286427	20.000	ng	-0.01
86) Perylene-d12	23.727	264	1236135	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.416	112	2085025	121.312	ng	0.00
7) Phenol-d6	7.011	99	2543785	122.222	ng	0.00
23) Nitrobenzene-d5	8.987	82	1570271	86.681	ng	-0.01
42) 2,4,6-Tribromophenol	15.969	330	1250118	125.543	ng	0.00
45) 2-Fluorobiphenyl	13.093	172	4169703	80.380	ng	0.00
79) Terphenyl-d14	19.792	244	6713184	88.085	ng	0.00

Target Compounds	Qvalue

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

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