

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM072122\
 Data File : BM035891.D
 Acq On : 21 Jul 2022 16:08
 Operator : CG/JU
 Sample : PB146352BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB146352BL

Quant Time: Jul 22 01:08:14 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM071822.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jul 18 17:08:15 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.898	152	311298	20.000	ng	0.00
21) Naphthalene-d8	10.704	136	1267969	20.000	ng	0.00
39) Acenaphthene-d10	14.527	164	677793	20.000	ng	0.00
64) Phenanthrene-d10	17.268	188	1270960	20.000	ng	0.00
76) Chrysene-d12	21.444	240	966668	20.000	ng	0.00
86) Perylene-d12	23.821	264	848871	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.457	112	2844075	141.638	ng	0.00
7) Phenol-d6	7.063	99	3436155	136.085	ng	0.00
23) Nitrobenzene-d5	9.063	82	2043929	86.214	ng	0.00
42) 2,4,6-Tribromophenol	16.015	330	971875	128.191	ng	0.00
45) 2-Fluorobiphenyl	13.163	172	4266207	88.970	ng	0.00
79) Terphenyl-d14	19.892	244	5110197	102.855	ng	0.00

Target Compounds	Qvalue

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

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM072122\
Data File : BM035891.D
Acq On : 21 Jul 2022 16:08
Operator : CG/JU
Sample : PB146352BL
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
PB146352BL

Quant Time: Jul 22 01:08:14 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM071822.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Jul 18 17:08:15 2022
Response via : Initial Calibration

