

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061021\
 Data File : BM030367.D
 Acq On : 10 Jun 2021 10:29
 Operator : CG/JU
 Sample : PB136969BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB136969BL

Quant Time: Jun 10 11:04:40 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM060921.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 08 20:51:08 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.845	152	212587	20.000	ng	0.00
21) Naphthalene-d8	10.633	136	846419	20.000	ng	0.00
39) Acenaphthene-d10	14.468	164	593519	20.000	ng	0.00
64) Phenanthrene-d10	17.204	188	1378990	20.000	ng	0.00
76) Chrysene-d12	21.368	240	1582853	20.000	ng	0.00
86) Perylene-d12	23.644	264	1671023	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.428	112	1884109	144.078	ng	0.00
7) Phenol-d6	7.010	99	2416275	127.638	ng	0.00
23) Nitrobenzene-d5	8.998	82	1380594	79.463	ng	-0.01
42) 2,4,6-Tribromophenol	15.951	330	1103251	147.846	ng	0.00
45) 2-Fluorobiphenyl	13.092	172	3653901	80.960	ng	0.00
79) Terphenyl-d14	19.821	244	7151680	79.360	ng	0.00

Target Compounds	Qvalue

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

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