

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP060921\
 Data File : BP005819.D
 Acq On : 09 Jun 2021 20:53
 Operator : CG/JU
 Sample : M2589-34
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 18-B11(0-2)

Quant Time: Jun 10 01:22:23 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP060821.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jun 09 11:54:55 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.952	152	71570	20.000	ng	0.00
21) Naphthalene-d8	10.757	136	265190	20.000	ng	0.00
39) Acenaphthene-d10	14.581	164	163266	20.000	ng	0.00
64) Phenanthrene-d10	17.322	188	344250	20.000	ng	0.00
76) Chrysene-d12	21.386	240	409336	20.000	ng	0.00
86) Perylene-d12	23.810	264	471818	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.516	112	534244	119.780	ng	0.00
7) Phenol-d6	7.116	99	570523	98.688	ng	0.00
23) Nitrobenzene-d5	9.116	82	444054	82.095	ng	0.00
42) 2,4,6-Tribromophenol	16.063	330	365729	130.535	ng	0.00
45) 2-Fluorobiphenyl	13.210	172	976134	78.509	ng	0.00
79) Terphenyl-d14	19.869	244	1807530	82.685	ng	0.00

Target Compounds	Qvalue

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

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