

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM020922\
 Data File : BM034034.D
 Acq On : 09 Feb 2022 19:32
 Operator : CG/JU
 Sample : SP5811
 Misc : 625 SURROGATE
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SP5811

Quant Time: Feb 10 04:07:16 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM020922.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Feb 09 17:52:36 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.900	152	153887	20.000	ng	0.00
21) Naphthalene-d8	10.715	136	561847	20.000	ng	# 0.00
39) Acenaphthene-d10	14.545	164	324315	20.000	ng	0.00
64) Phenanthrene-d10	17.270	188	624786	20.000	ng	0.00
76) Chrysene-d12	21.437	240	457524	20.000	ng	# 0.00
86) Perylene-d12	23.825	264	425645	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.444	112	1042192	109.659	ng	0.00
7) Phenol-d6	7.066	99	1276414	110.169	ng	0.00
23) Nitrobenzene-d5	9.071	82	1210553	101.205	ng	0.00
42) 2,4,6-Tribromophenol	16.031	330	372387	103.727	ng	0.00
45) 2-Fluorobiphenyl	13.171	172	2781571	112.607	ng	0.00
79) Terphenyl-d14	19.883	244	2445462	105.711	ng	0.00

Target Compounds	Qvalue

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

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