

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM042722\
 Data File : BM034861.D
 Acq On : 27 Apr 2022 20:35
 Operator : CG/JU
 Sample : SP5883
 Misc : 8270 SURROGATE
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SP5883

Quant Time: Apr 28 02:01:08 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM042722.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Apr 28 01:51:59 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.934	152	227724	20.000	ng	0.00
21) Naphthalene-d8	10.733	136	887308	20.000	ng	0.00
39) Acenaphthene-d10	14.563	164	525932	20.000	ng	0.00
64) Phenanthrene-d10	17.304	188	1030058	20.000	ng	0.00
76) Chrysene-d12	21.480	240	729100	20.000	ng	0.00
86) Perylene-d12	23.868	264	620642	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.510	112	2238686	173.435	ng	0.00
7) Phenol-d6	7.098	99	2771932	166.349	ng	0.00
23) Nitrobenzene-d5	9.092	82	1788205	108.830	ng	0.00
42) 2,4,6-Tribromophenol	16.051	330	952242	166.444	ng	0.00
45) 2-Fluorobiphenyl	13.192	172	4013857	110.790	ng	0.00
79) Terphenyl-d14	19.927	244	4961851	131.768	ng	0.00

Target Compounds	Qvalue

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

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