

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020122\
 Data File : BP008989.D
 Acq On : 01 Feb 2022 16:30
 Operator : CG/JU
 Sample : N1335-07
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 MH-4S

Quant Time: Feb 01 17:25:36 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP011422.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Feb 01 04:16:03 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.828	152	234501	20.000	ng	0.00
21) Naphthalene-d8	10.628	136	910441	20.000	ng	0.00
39) Acenaphthene-d10	14.469	164	585030	20.000	ng	0.00
64) Phenanthrene-d10	17.222	188	1191407	20.000	ng	-0.01
76) Chrysene-d12	21.316	240	1103580	20.000	ng	-0.01
86) Perylene-d12	23.721	264	1210983	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.417	112	1474080	97.208	ng	0.00
7) Phenol-d6	7.011	99	1812856	98.724	ng	0.00
23) Nitrobenzene-d5	8.987	82	1119516	69.811	ng	-0.01
42) 2,4,6-Tribromophenol	15.963	330	835083	98.064	ng	-0.01
45) 2-Fluorobiphenyl	13.087	172	2834931	63.903	ng	-0.01
79) Terphenyl-d14	19.786	244	4034111	61.702	ng	-0.01
Target Compounds						
88) Benzo(b)fluoranthene	22.992	252	182823	2.261	ng	Qvalue # 94

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

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