

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110321\
 Data File : BP007832.D
 Acq On : 03 Nov 2021 20:27
 Operator : CG/JU
 Sample : M4395-07
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 MH-GGS

Quant Time: Nov 04 07:31:55 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP110221.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 02 19:18:55 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.916	152	215125	20.00	ng	0.00
21) Naphthalene-d8	10.722	136	879698	20.00	ng	0.00
39) Acenaphthene-d10	14.557	164	565595	20.00	ng	0.00
64) Phenanthrene-d10	17.310	188	1229006	20.00	ng	-0.01
76) Chrysene-d12	21.404	240	1377722	20.00	ng	-0.01
86) Perylene-d12	23.839	264	1271091	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.499	112	1559404	122.94	ng	0.00
7) Phenol-d6	7.093	99	2004394	109.17	ng	0.00
23) Nitrobenzene-d5	9.075	82	1220485	74.53	ng	-0.01
42) 2,4,6-Tribromophenol	16.051	330	906922	112.36	ng	0.00
45) 2-Fluorobiphenyl	13.181	172	2727209	70.22	ng	-0.01
79) Terphenyl-d14	19.875	244	4591277	67.86	ng	0.00
Target Compounds						
						Qvalue
75) Fluoranthene	19.322	202	252851	3.10	ng	98
78) Pyrene	19.675	202	217232	2.57	ng	100
88) Benzo(b)fluoranthene	23.098	252	175807	2.35	ng	97

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

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