

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111121\
 Data File : BP007946.D
 Acq On : 11 Nov 2021 15:32
 Operator : CG/JU
 Sample : M4607-03
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 11966

Quant Time: Nov 11 17:39:53 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP110221.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Nov 11 13:53:53 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.846	152	157377	20.000	ng	0.00
21) Naphthalene-d8	10.646	136	543073	20.000	ng	0.00
39) Acenaphthene-d10	14.487	164	375363	20.000	ng	0.00
64) Phenanthrene-d10	17.239	188	912406	20.000	ng	# 0.00
76) Chrysene-d12	21.333	240	1223409	20.000	ng	# 0.00
86) Perylene-d12	23.739	264	1285809	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.434	112	437220	47.116	ng	0.00
7) Phenol-d6	7.028	99	1074667	80.012	ng	0.00
23) Nitrobenzene-d5	9.004	82	805694	79.694	ng	0.00
42) 2,4,6-Tribromophenol	15.981	330	102693	19.170	ng	0.00
45) 2-Fluorobiphenyl	13.110	172	2018826	78.323	ng	0.00
79) Terphenyl-d14	19.810	244	4487558	74.693	ng	0.00

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

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