

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP120721\
 Data File : BP008248.D
 Acq On : 07 Dec 2021 13:08
 Operator : CG/JU
 Sample : M4919-07
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 EB-20211202

Quant Time: Dec 08 01:07:46 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP112521.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 03 14:39:24 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.787	152	31154	20.000	ng	-0.01
21) Naphthalene-d8	10.581	136	114014	20.000	ng	-0.02
39) Acenaphthene-d10	14.434	164	75102	20.000	ng	-0.01
64) Phenanthrene-d10	17.198	188	173681	20.000	ng	-0.01
76) Chrysene-d12	21.304	240	235057	20.000	ng	-0.01
86) Perylene-d12	23.698	264	291729	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.381	112	139216	71.949	ng	0.00
7) Phenol-d6	6.975	99	107150	41.022	ng	-0.01
23) Nitrobenzene-d5	8.946	82	267227	106.580	ng	-0.02
42) 2,4,6-Tribromophenol	15.939	330	167021	173.993	ng	-0.01
45) 2-Fluorobiphenyl	13.051	172	561300	108.560	ng	-0.02
79) Terphenyl-d14	19.769	244	1178829	104.811	ng	-0.01

Target Compounds	Qvalue

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

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