

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP120721\
 Data File : BP008249.D
 Acq On : 07 Dec 2021 13:42
 Operator : CG/JU
 Sample : M4927-07
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PH1-BOT-9

Quant Time: Dec 08 01:07:58 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	21238	20.000	ng	-0.01
21) Naphthalene-d8	10.587	136	83056	20.000	ng	-0.01
39) Acenaphthene-d10	14.439	164	53412	20.000	ng	0.00
64) Phenanthrene-d10	17.198	188	121265	20.000	ng	-0.01
76) Chrysene-d12	21.304	240	182632	20.000	ng	-0.01
86) Perylene-d12	23.704	264	231201	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.381	112	157825	119.650	ng	0.00
7) Phenol-d6	6.975	99	216465	121.567	ng	-0.01
23) Nitrobenzene-d5	8.946	82	140331	76.831	ng	-0.02
42) 2,4,6-Tribromophenol	15.939	330	81805	119.826	ng	-0.01
45) 2-Fluorobiphenyl	13.051	172	286644	77.953	ng	-0.02
79) Terphenyl-d14	19.769	244	647801	74.130	ng	-0.01
Target Compounds						
10) Phenol	6.999	94	4220	2.306	ng	Qvalue # 49

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

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