

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111021\
 Data File : BP007922.D
 Acq On : 10 Nov 2021 16:31
 Operator : CG/JU
 Sample : PB140662BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB140662BL

Quant Time: Nov 11 10:14:15 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 09 10:45:24 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.858	152	132973	20.000	ng	0.00
21) Naphthalene-d8	10.657	136	494754	20.000	ng	# 0.00
39) Acenaphthene-d10	14.498	164	373429	20.000	ng	0.00
64) Phenanthrene-d10	17.251	188	924300	20.000	ng	# 0.00
76) Chrysene-d12	21.345	240	1170956	20.000	ng	0.00
86) Perylene-d12	23.757	264	1230010	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.446	112	1073155	136.870	ng	0.00
7) Phenol-d6	7.040	99	1326325	116.871	ng	0.00
23) Nitrobenzene-d5	9.016	82	752868	81.742	ng	-0.01
42) 2,4,6-Tribromophenol	15.992	330	777566	145.902	ng	0.00
45) 2-Fluorobiphenyl	13.122	172	2046351	79.802	ng	0.00
79) Terphenyl-d14	19.816	244	4532458	78.819	ng	0.00

Target Compounds	Qvalue

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

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