

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080221\
 Data File : BP006457.D
 Acq On : 02 Aug 2021 18:56
 Operator : CG/JU
 Sample : PB138103BL
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB138103BL

Quant Time: Aug 03 05:07:49 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP080221.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Aug 03 05:06:41 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.758	152	235465	20.000	ng	0.00
21) Naphthalene-d8	10.546	136	1017977	20.000	ng	# 0.00
39) Acenaphthene-d10	14.387	164	565067	20.000	ng	0.00
64) Phenanthrene-d10	17.140	188	1143105	20.000	ng	# 0.00
76) Chrysene-d12	21.239	240	1031349	20.000	ng	# 0.00
86) Perylene-d12	23.569	264	1037612	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.358	112	1773972	117.489	ng	0.00
7) Phenol-d6	6.940	99	2263353	106.176	ng	0.00
23) Nitrobenzene-d5	8.916	82	1469269	68.047	ng	0.00
42) 2,4,6-Tribromophenol	15.881	330	602575	104.289	ng	0.00
45) 2-Fluorobiphenyl	13.016	172	2567354	68.102	ng	0.00
79) Terphenyl-d14	19.716	244	3685034	71.946	ng	0.00

Target Compounds	Qvalue

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

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