

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080421\
 Data File : BP006496.D
 Acq On : 04 Aug 2021 17:13
 Operator : CG/JU
 Sample : PB138118BL
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB138118BL

Quant Time: Aug 04 17:46:03 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP080321.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Aug 03 15:41:30 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.752	152	189084	20.000	ng	0.00
21) Naphthalene-d8	10.534	136	794020	20.000	ng	#-0.01
39) Acenaphthene-d10	14.387	164	426585	20.000	ng	0.00
64) Phenanthrene-d10	17.139	188	833647	20.000	ng	# 0.00
76) Chrysene-d12	21.239	240	736407	20.000	ng	# 0.00
86) Perylene-d12	23.568	264	734411	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.352	112	1820044	147.843	ng	0.00
7) Phenol-d6	6.928	99	2341977	133.923	ng	0.00
23) Nitrobenzene-d5	8.910	82	1544050	89.758	ng	0.00
42) 2,4,6-Tribromophenol	15.881	330	542669	121.722	ng	0.00
45) 2-Fluorobiphenyl	13.010	172	2590926	90.869	ng	0.00
79) Terphenyl-d14	19.716	244	3389921	92.241	ng	0.00

Target Compounds	Qvalue

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

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