

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082421\
 Data File : BP006861.D
 Acq On : 24 Aug 2021 23:22
 Operator : CG/JU
 Sample : M3489-16
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 1-GROWS

Quant Time: Aug 25 02:36:42 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 24 13:40:38 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.687	152	166350	20.000	ng	0.00
21) Naphthalene-d8	10.469	136	677472	20.000	ng	# 0.00
39) Acenaphthene-d10	14.328	164	384375	20.000	ng	-0.01
64) Phenanthrene-d10	17.086	188	765549	20.000	ng	#-0.01
76) Chrysene-d12	21.192	240	759697	20.000	ng	#-0.02
86) Perylene-d12	23.498	264	811604	20.000	ng	-0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.293	112	1273488	117.584	ng	0.00
7) Phenol-d6	6.881	99	1434110	93.215	ng	0.00
23) Nitrobenzene-d5	8.846	82	1183889	80.661	ng	0.00
42) 2,4,6-Tribromophenol	15.828	330	600383	149.455	ng	-0.02
45) 2-Fluorobiphenyl	12.951	172	2181554	84.914	ng	0.00
79) Terphenyl-d14	19.669	244	3395881	89.570	ng	-0.02
Target Compounds						
32) Benzoic acid	9.775	122	51235	7.123	ng	Qvalue 87

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

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