

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081721\
 Data File : BP006696.D
 Acq On : 16 Aug 2021 22:37
 Operator : CG/JU
 Sample : M3424-09
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 19N

Quant Time: Aug 17 06:44:22 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP080321.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 16 18:17:59 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.705	152	178441	20.000	ng	0.00
21) Naphthalene-d8	10.487	136	741246	20.000	ng	# 0.00
39) Acenaphthene-d10	14.345	164	402169	20.000	ng	0.00
64) Phenanthrene-d10	17.104	188	805952	20.000	ng	# 0.00
76) Chrysene-d12	21.204	240	801331	20.000	ng	# 0.00
86) Perylene-d12	23.516	264	853726	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.311	112	1610844	138.654	ng	0.00
7) Phenol-d6	6.893	99	1888319	114.422	ng	0.00
23) Nitrobenzene-d5	8.863	82	1486919	92.591	ng	0.00
42) 2,4,6-Tribromophenol	15.845	330	632876	150.573	ng	0.00
45) 2-Fluorobiphenyl	12.969	172	2442193	90.853	ng	0.00
79) Terphenyl-d14	19.686	244	3916480	97.934	ng	0.00

Target Compounds	Qvalue

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

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