

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041321\
 Data File : BP005285.D
 Acq On : 13 Apr 2021 14:02
 Operator : CG/JU
 Sample : M1967-13RE
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 EB 2021-04-06 SEDRE

Quant Time: Apr 14 02:41:26 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP040921.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Apr 09 18:44:11 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.675	152	40049	20.00	ng	# 0.00
21) Naphthalene-d8	10.463	136	129581	20.00	ng	# 0.00
39) Acenaphthene-d10	14.340	164	81785	20.00	ng	0.00
64) Phenanthrene-d10	17.098	188	200721	20.00	ng	# 0.00
76) Chrysene-d12	21.204	240	269161	20.00	ng	# 0.00
86) Perylene-d12	23.474	264	275977	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.275	112	226286	59.30	ng	0.00
7) Phenol-d6	6.864	99	137576	31.60	ng	0.00
23) Nitrobenzene-d5	8.840	82	395008	96.13	ng	0.00
42) 2,4,6-Tribromophenol	15.845	330	173935	168.48	ng	0.00
45) 2-Fluorobiphenyl	12.951	172	607287	100.25	ng	0.00
79) Terphenyl-d14	19.681	244	1297366	94.69	ng	0.00

Target Compounds	Qvalue

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

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