

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040121\
 Data File : BP005134.D
 Acq On : 01 Apr 2021 14:02
 Operator : CG/JU
 Sample : M1864-01
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 MMR-SP

Quant Time: Apr 01 14:43:21 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP032321.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Apr 01 12:23:57 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.705	152	26292	20.00	ng	# 0.00
21) Naphthalene-d8	10.493	136	97517	20.00	ng	# 0.00
39) Acenaphthene-d10	14.363	164	61825	20.00	ng	0.00
64) Phenanthrene-d10	17.122	188	136985	20.00	ng	# 0.00
76) Chrysene-d12	21.222	240	147275	20.00	ng	0.00
86) Perylene-d12	23.504	264	155257	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.305	112	227597	133.22	ng	0.00
7) Phenol-d6	6.887	99	264675	108.15	ng	0.00
23) Nitrobenzene-d5	8.863	82	210811	92.76	ng	0.00
42) 2,4,6-Tribromophenol	15.863	330	117733	146.16	ng	0.00
45) 2-Fluorobiphenyl	12.975	172	416269	91.08	ng	0.00
79) Terphenyl-d14	19.698	244	749857	92.69	ng	0.00

Target Compounds	Qvalue

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

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