Quantitation Report (QT Reviewed)

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020122\
Data File : BP008984.D
Acq On : 01 Feb 2022 13:44
Operator : CG/JU
Sample : N1303-07RE
Misc :
ALS Vial : 7 Sample Multiplier: 1

Quant Time: Feb 01 15:30:32 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP011422.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Feb 01 04:16:03 2022
Response via : Initial Calibration

Instrument :
BNA_P
ClientSampleId :
CC-8RE

Manual Integrations APPROVED

Reviewed By :Christian Giraldo 02/02/2022
Supervised By :Jagrut Upadhyay 02/02/2022

Compound	R.T.	QIon	Response	Conc	Units		Dev(Min)
Internal Standards							
1) 1,4-Dichlorobenzene-d4	7.828	152	217436	20.000	ng		0.00
21) Naphthalene-d8	10.622	136	874561	20.000	ng		-0.01
39) Acenaphthene-d10	14.469	164	570485	20.000	ng		0.00
64) Phenanthrene-d10	17.222	188	1176610	20.000	ng		-0.01
76) Chrysene-d12	21.316	240	1137364	20.000	ng		-0.01
86) Perylene-d12	23.721	264	1181282	20.000	ng		-0.02
System Monitoring Compounds							
5) 2-Fluorophenol	5.422	112	1631	0.116	ng		0.00
7) Phenol-d6	7.016	99	48190	2.830	ng		0.00
23) Nitrobenzene-d5	8.987	82	1206454	78.319	ng		-0.01
42) 2,4,6-Tribromophenol	15.975	330	1081	0.130	ng		0.00
45) 2-Fluorobiphenyl	13.087	172	3114568	71.997	ng		-0.01
79) Terphenyl-d14	19.786	244	4746192	70.437	ng		-0.01
Target Compounds							Qvalue
31) Naphthalene	10.675	128	1603368	33.755	ng		99
37) 2-Methylnaphthalene	12.287	142	202666	5.836	ng		98
38) 1-Methylnaphthalene	12.504	142	111358m	3.494	ng		
58) Fluorene	15.516	166	88376	2.141	ng		97
71) Phenanthrene	17.263	178	975643	14.727	ng		99
72) Anthracene	17.357	178	172621	2.597	ng		99
75) Fluoranthene	19.233	202	1049348	14.244	ng		97
78) Pyrene	19.586	202	980779	12.359	ng		98
81) Benzo(a)anthracene	21.298	228	585923	7.657	ng		98
83) Chrysene	21.351	228	605570	8.164	ng		98
87) Indeno(1,2,3-cd)pyrene	26.221	276	321210	3.300	ng	#	93
88) Benzo(b)fluoranthene	22.986	252	665595	8.437	ng		97
89) Benzo(k)fluoranthene	23.027	252	299118m	3.917	ng		
90) Benzo(a)pyrene	23.616	252	464184	6.096	ng		98
92) Benzo(g,h,i)perylene	26.992	276	309683	3.832	ng		97

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

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020122\
Data File : BP008984.D
Acq On : 01 Feb 2022 13:44
Operator : CG/JU
Sample : N1303-07RE
Misc :
ALS Vial : 7 Sample Multiplier: 1

Quant Time: Feb 01 15:30:32 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP011422.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Feb 01 04:16:03 2022
Response via : Initial Calibration

Instrument :
BNA_P
ClientSampleId :
CC-8RE

Manual Integrations APPROVED

Reviewed By :Christian Giraldo 02/02/2022
Supervised By :Jagrut Upadhyay 02/02/2022